



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:58 AM GMT

PDB ID : 3GMH
Title : Crystal Structure of the Mad2 Dimer
Authors : Ozkan, E.; Luo, X.; Machius, M.; Yu, H.; Deisenhofer, J.
Deposited on : 2009-03-13
Resolution : 3.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

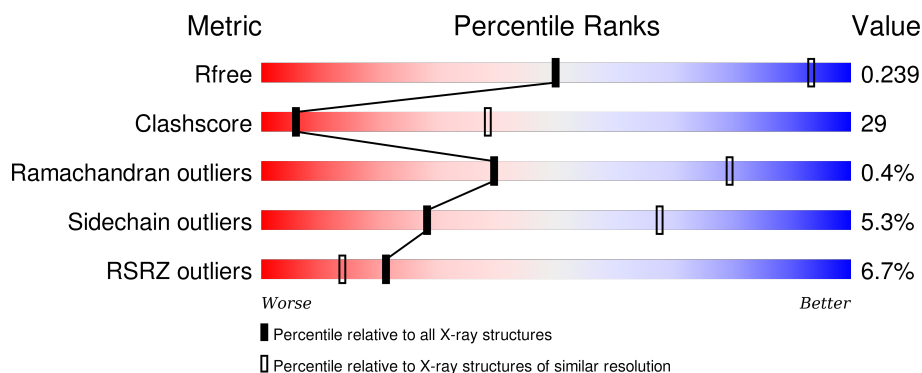
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1011 (4.38-3.54)
Clashscore	102246	1025 (4.32-3.60)
Ramachandran outliers	100387	1008 (4.34-3.58)
Sidechain outliers	100360	1027 (4.36-3.56)
RSRZ outliers	91569	1015 (4.38-3.54)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	207	<div> <div>2%</div> <div> <div></div> <div>56%</div> <div>35%</div> <div>7%</div> </div> </div>
1	B	207	<div> <div>0%</div> <div> <div></div> <div>44%</div> <div>41%</div> <div>11%</div> </div> </div>
1	C	207	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>36%</div> <div>6%</div> </div> </div>
1	D	207	<div> <div>2%</div> <div> <div></div> <div>47%</div> <div>39%</div> <div>10%</div> </div> </div>
1	E	207	<div> <div>9%</div> <div> <div></div> <div>55%</div> <div>34%</div> <div>9%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	207	
1	G	207	
1	H	207	
1	I	207	
1	J	207	
1	K	207	
1	L	207	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	207	-	-	X	-
2	SO4	C	206	-	-	X	-
2	SO4	D	206	-	-	-	X
2	SO4	E	206	-	-	X	-
2	SO4	G	206	-	-	-	X
2	SO4	K	206	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18301 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Mitotic spindle assembly checkpoint protein MAD2A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1561	1003	253	301	4			
1	B	185	Total	C	N	O	S	0	0	0
			1495	965	240	286	4			
1	C	195	Total	C	N	O	S	0	0	0
			1576	1012	256	304	4			
1	D	187	Total	C	N	O	S	0	0	0
			1515	977	246	288	4			
1	E	188	Total	C	N	O	S	0	0	0
			1522	981	245	292	4			
1	F	185	Total	C	N	O	S	0	0	0
			1495	965	240	286	4			
1	G	187	Total	C	N	O	S	0	0	0
			1518	978	246	290	4			
1	H	186	Total	C	N	O	S	0	0	0
			1505	971	243	287	4			
1	I	188	Total	C	N	O	S	0	0	0
			1525	983	247	291	4			
1	J	185	Total	C	N	O	S	0	0	0
			1495	965	240	286	4			
1	K	188	Total	C	N	O	S	0	0	0
			1525	983	247	291	4			
1	L	187	Total	C	N	O	S	0	0	0
			1509	974	243	288	4			

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP Q13257
A	0	ARG	-	EXPRESSION TAG	UNP Q13257
A	1	GLY	-	EXPRESSION TAG	UNP Q13257
A	2	SER	-	EXPRESSION TAG	UNP Q13257
A	3	HIS	-	EXPRESSION TAG	UNP Q13257

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Chain	Residue	Modelled	Actual	Comment	Reference
A	4	HIS	-	EXPRESSION TAG	UNP Q13257
A	5	HIS	-	EXPRESSION TAG	UNP Q13257
A	6	HIS	-	EXPRESSION TAG	UNP Q13257
A	7	HIS	-	EXPRESSION TAG	UNP Q13257
A	8	HIS	-	EXPRESSION TAG	UNP Q13257
A	9	GLY	-	EXPRESSION TAG	UNP Q13257
A	10	SER	-	EXPRESSION TAG	UNP Q13257
B	-1	MET	-	EXPRESSION TAG	UNP Q13257
B	0	ARG	-	EXPRESSION TAG	UNP Q13257
B	1	GLY	-	EXPRESSION TAG	UNP Q13257
B	2	SER	-	EXPRESSION TAG	UNP Q13257
B	3	HIS	-	EXPRESSION TAG	UNP Q13257
B	4	HIS	-	EXPRESSION TAG	UNP Q13257
B	5	HIS	-	EXPRESSION TAG	UNP Q13257
B	6	HIS	-	EXPRESSION TAG	UNP Q13257
B	7	HIS	-	EXPRESSION TAG	UNP Q13257
B	8	HIS	-	EXPRESSION TAG	UNP Q13257
B	9	GLY	-	EXPRESSION TAG	UNP Q13257
B	10	SER	-	EXPRESSION TAG	UNP Q13257
C	-1	MET	-	EXPRESSION TAG	UNP Q13257
C	0	ARG	-	EXPRESSION TAG	UNP Q13257
C	1	GLY	-	EXPRESSION TAG	UNP Q13257
C	2	SER	-	EXPRESSION TAG	UNP Q13257
C	3	HIS	-	EXPRESSION TAG	UNP Q13257
C	4	HIS	-	EXPRESSION TAG	UNP Q13257
C	5	HIS	-	EXPRESSION TAG	UNP Q13257
C	6	HIS	-	EXPRESSION TAG	UNP Q13257
C	7	HIS	-	EXPRESSION TAG	UNP Q13257
C	8	HIS	-	EXPRESSION TAG	UNP Q13257
C	9	GLY	-	EXPRESSION TAG	UNP Q13257
C	10	SER	-	EXPRESSION TAG	UNP Q13257
D	-1	MET	-	EXPRESSION TAG	UNP Q13257
D	0	ARG	-	EXPRESSION TAG	UNP Q13257
D	1	GLY	-	EXPRESSION TAG	UNP Q13257
D	2	SER	-	EXPRESSION TAG	UNP Q13257
D	3	HIS	-	EXPRESSION TAG	UNP Q13257
D	4	HIS	-	EXPRESSION TAG	UNP Q13257
D	5	HIS	-	EXPRESSION TAG	UNP Q13257
D	6	HIS	-	EXPRESSION TAG	UNP Q13257
D	7	HIS	-	EXPRESSION TAG	UNP Q13257
D	8	HIS	-	EXPRESSION TAG	UNP Q13257
D	9	GLY	-	EXPRESSION TAG	UNP Q13257

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Chain	Residue	Modelled	Actual	Comment	Reference
D	10	SER	-	EXPRESSION TAG	UNP Q13257
E	-1	MET	-	EXPRESSION TAG	UNP Q13257
E	0	ARG	-	EXPRESSION TAG	UNP Q13257
E	1	GLY	-	EXPRESSION TAG	UNP Q13257
E	2	SER	-	EXPRESSION TAG	UNP Q13257
E	3	HIS	-	EXPRESSION TAG	UNP Q13257
E	4	HIS	-	EXPRESSION TAG	UNP Q13257
E	5	HIS	-	EXPRESSION TAG	UNP Q13257
E	6	HIS	-	EXPRESSION TAG	UNP Q13257
E	7	HIS	-	EXPRESSION TAG	UNP Q13257
E	8	HIS	-	EXPRESSION TAG	UNP Q13257
E	9	GLY	-	EXPRESSION TAG	UNP Q13257
E	10	SER	-	EXPRESSION TAG	UNP Q13257
F	-1	MET	-	EXPRESSION TAG	UNP Q13257
F	0	ARG	-	EXPRESSION TAG	UNP Q13257
F	1	GLY	-	EXPRESSION TAG	UNP Q13257
F	2	SER	-	EXPRESSION TAG	UNP Q13257
F	3	HIS	-	EXPRESSION TAG	UNP Q13257
F	4	HIS	-	EXPRESSION TAG	UNP Q13257
F	5	HIS	-	EXPRESSION TAG	UNP Q13257
F	6	HIS	-	EXPRESSION TAG	UNP Q13257
F	7	HIS	-	EXPRESSION TAG	UNP Q13257
F	8	HIS	-	EXPRESSION TAG	UNP Q13257
F	9	GLY	-	EXPRESSION TAG	UNP Q13257
F	10	SER	-	EXPRESSION TAG	UNP Q13257
G	-1	MET	-	EXPRESSION TAG	UNP Q13257
G	0	ARG	-	EXPRESSION TAG	UNP Q13257
G	1	GLY	-	EXPRESSION TAG	UNP Q13257
G	2	SER	-	EXPRESSION TAG	UNP Q13257
G	3	HIS	-	EXPRESSION TAG	UNP Q13257
G	4	HIS	-	EXPRESSION TAG	UNP Q13257
G	5	HIS	-	EXPRESSION TAG	UNP Q13257
G	6	HIS	-	EXPRESSION TAG	UNP Q13257
G	7	HIS	-	EXPRESSION TAG	UNP Q13257
G	8	HIS	-	EXPRESSION TAG	UNP Q13257
G	9	GLY	-	EXPRESSION TAG	UNP Q13257
G	10	SER	-	EXPRESSION TAG	UNP Q13257
H	-1	MET	-	EXPRESSION TAG	UNP Q13257
H	0	ARG	-	EXPRESSION TAG	UNP Q13257
H	1	GLY	-	EXPRESSION TAG	UNP Q13257
H	2	SER	-	EXPRESSION TAG	UNP Q13257
H	3	HIS	-	EXPRESSION TAG	UNP Q13257

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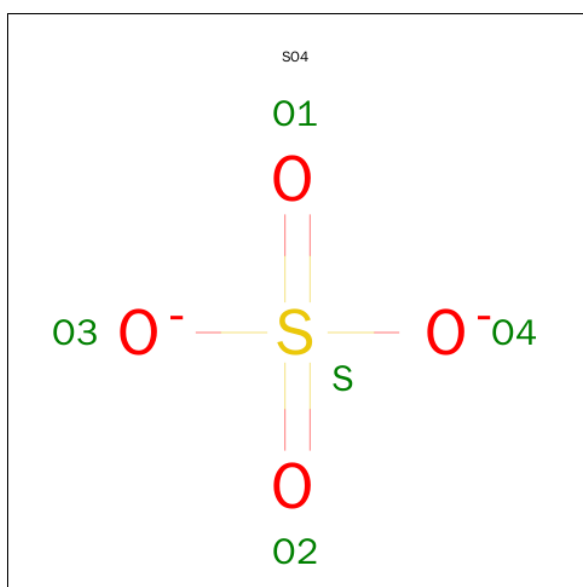
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H	4	HIS	-	EXPRESSION TAG	UNP Q13257
H	5	HIS	-	EXPRESSION TAG	UNP Q13257
H	6	HIS	-	EXPRESSION TAG	UNP Q13257
H	7	HIS	-	EXPRESSION TAG	UNP Q13257
H	8	HIS	-	EXPRESSION TAG	UNP Q13257
H	9	GLY	-	EXPRESSION TAG	UNP Q13257
H	10	SER	-	EXPRESSION TAG	UNP Q13257
I	-1	MET	-	EXPRESSION TAG	UNP Q13257
I	0	ARG	-	EXPRESSION TAG	UNP Q13257
I	1	GLY	-	EXPRESSION TAG	UNP Q13257
I	2	SER	-	EXPRESSION TAG	UNP Q13257
I	3	HIS	-	EXPRESSION TAG	UNP Q13257
I	4	HIS	-	EXPRESSION TAG	UNP Q13257
I	5	HIS	-	EXPRESSION TAG	UNP Q13257
I	6	HIS	-	EXPRESSION TAG	UNP Q13257
I	7	HIS	-	EXPRESSION TAG	UNP Q13257
I	8	HIS	-	EXPRESSION TAG	UNP Q13257
I	9	GLY	-	EXPRESSION TAG	UNP Q13257
I	10	SER	-	EXPRESSION TAG	UNP Q13257
J	-1	MET	-	EXPRESSION TAG	UNP Q13257
J	0	ARG	-	EXPRESSION TAG	UNP Q13257
J	1	GLY	-	EXPRESSION TAG	UNP Q13257
J	2	SER	-	EXPRESSION TAG	UNP Q13257
J	3	HIS	-	EXPRESSION TAG	UNP Q13257
J	4	HIS	-	EXPRESSION TAG	UNP Q13257
J	5	HIS	-	EXPRESSION TAG	UNP Q13257
J	6	HIS	-	EXPRESSION TAG	UNP Q13257
J	7	HIS	-	EXPRESSION TAG	UNP Q13257
J	8	HIS	-	EXPRESSION TAG	UNP Q13257
J	9	GLY	-	EXPRESSION TAG	UNP Q13257
J	10	SER	-	EXPRESSION TAG	UNP Q13257
K	-1	MET	-	EXPRESSION TAG	UNP Q13257
K	0	ARG	-	EXPRESSION TAG	UNP Q13257
K	1	GLY	-	EXPRESSION TAG	UNP Q13257
K	2	SER	-	EXPRESSION TAG	UNP Q13257
K	3	HIS	-	EXPRESSION TAG	UNP Q13257
K	4	HIS	-	EXPRESSION TAG	UNP Q13257
K	5	HIS	-	EXPRESSION TAG	UNP Q13257
K	6	HIS	-	EXPRESSION TAG	UNP Q13257
K	7	HIS	-	EXPRESSION TAG	UNP Q13257
K	8	HIS	-	EXPRESSION TAG	UNP Q13257
K	9	GLY	-	EXPRESSION TAG	UNP Q13257

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Chain	Residue	Modelled	Actual	Comment	Reference
K	10	SER	-	EXPRESSION TAG	UNP Q13257
L	-1	MET	-	EXPRESSION TAG	UNP Q13257
L	0	ARG	-	EXPRESSION TAG	UNP Q13257
L	1	GLY	-	EXPRESSION TAG	UNP Q13257
L	2	SER	-	EXPRESSION TAG	UNP Q13257
L	3	HIS	-	EXPRESSION TAG	UNP Q13257
L	4	HIS	-	EXPRESSION TAG	UNP Q13257
L	5	HIS	-	EXPRESSION TAG	UNP Q13257
L	6	HIS	-	EXPRESSION TAG	UNP Q13257
L	7	HIS	-	EXPRESSION TAG	UNP Q13257
L	8	HIS	-	EXPRESSION TAG	UNP Q13257
L	9	GLY	-	EXPRESSION TAG	UNP Q13257
L	10	SER	-	EXPRESSION TAG	UNP Q13257

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	H	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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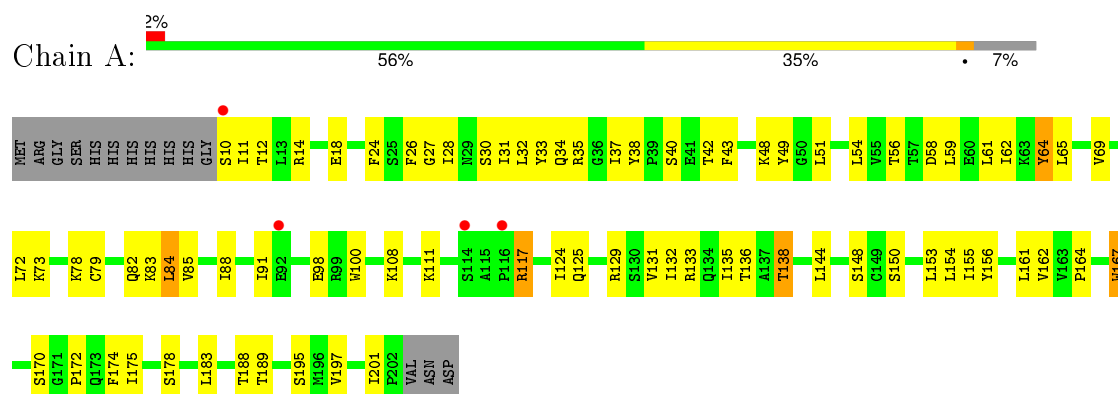
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		

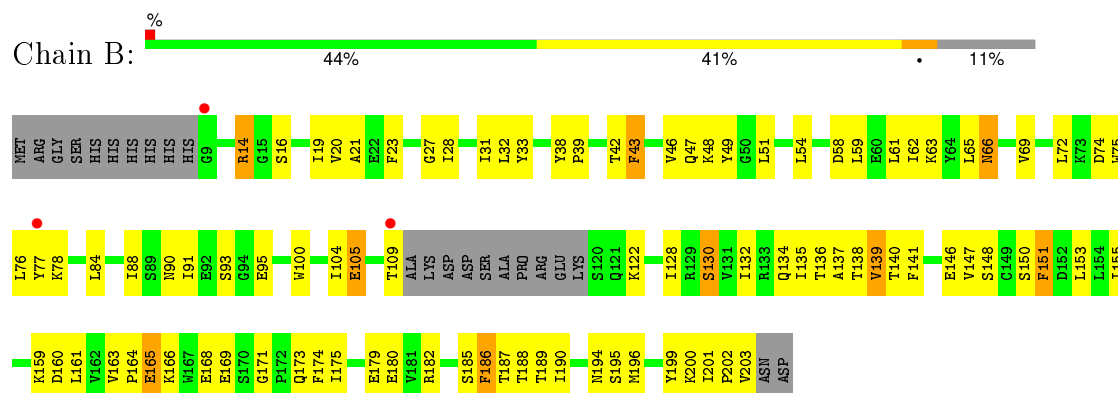
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

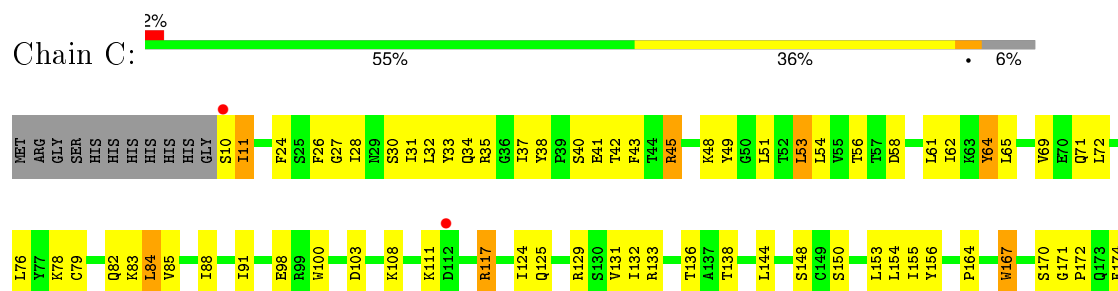
- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

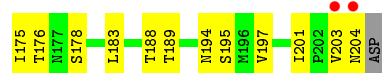


- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

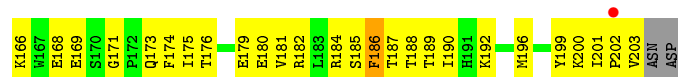


- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

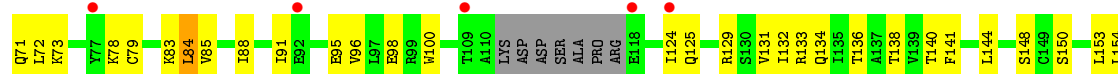
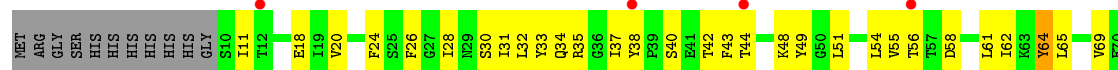




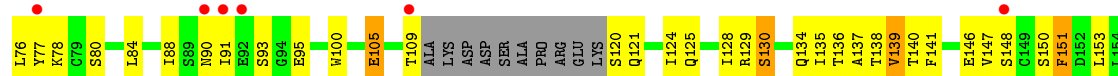
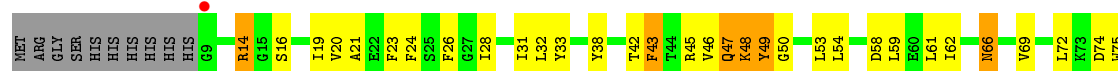
- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A



- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A



- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

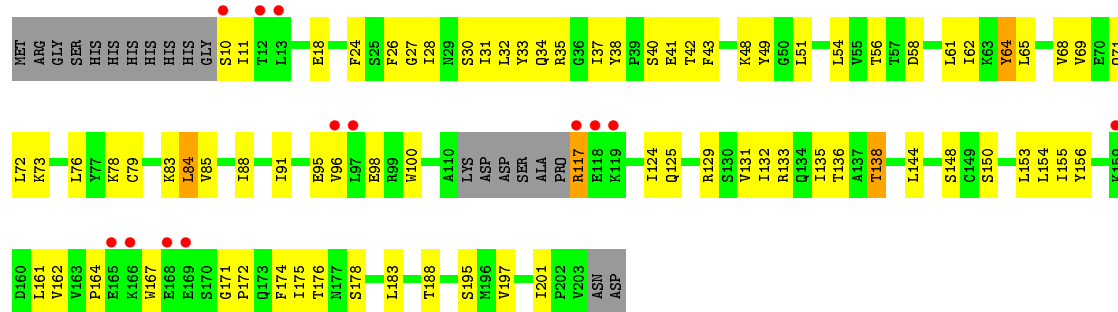


- Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A



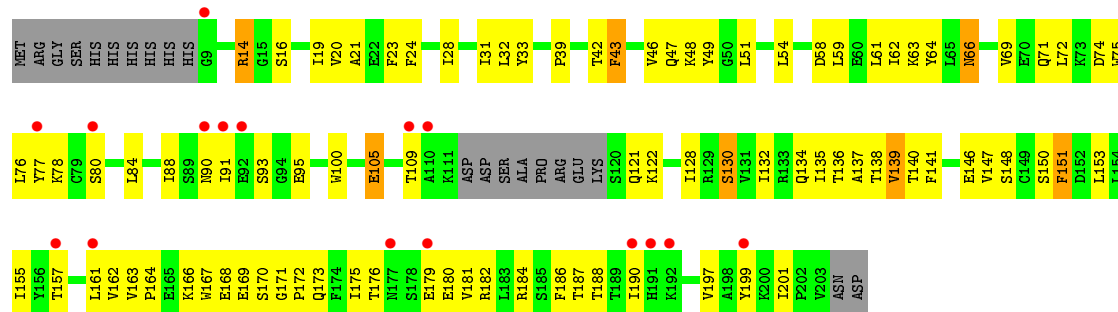
• Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

Chain K: 



• Molecule 1: Mitotic spindle assembly checkpoint protein MAD2A

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	124.83Å 104.92Å 157.75Å 90.00° 97.57° 90.00°	Depositor
Resolution (Å)	49.73 – 3.95 49.73 – 3.94	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.73-3.95) 98.8 (49.73-3.94)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 3.88Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 2009_02_15_2320_3)	Depositor
R, R_{free}	0.218 , 0.251 0.203 , 0.239	Depositor DCC
R_{free} test set	1769 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	115.9	Xtriage
Anisotropy	0.497	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 104.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 35706 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	18301	wwPDB-VP
Average B, all atoms (Å ²)	156.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.40	0/1589	0.55	0/2155
1	B	0.50	0/1521	0.61	0/2063
1	C	0.38	0/1604	0.54	0/2176
1	D	0.50	0/1543	0.61	0/2093
1	E	0.35	0/1548	0.52	0/2099
1	F	0.40	0/1521	0.57	0/2063
1	G	0.34	0/1544	0.52	0/2092
1	H	0.38	0/1532	0.55	0/2078
1	I	0.38	0/1551	0.55	0/2102
1	J	0.37	0/1521	0.56	0/2063
1	K	0.36	0/1551	0.53	0/2102
1	L	0.41	0/1535	0.57	0/2081
All	All	0.40	0/18560	0.56	0/25167

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1561	0	1581	87	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1495	0	1517	97	0
1	C	1576	0	1596	89	0
1	D	1515	0	1531	93	0
1	E	1522	0	1544	82	0
1	F	1495	0	1517	100	0
1	G	1518	0	1542	83	0
1	H	1505	0	1524	102	0
1	I	1525	0	1551	144	0
1	J	1495	0	1517	104	0
1	K	1525	0	1551	84	0
1	L	1509	0	1535	97	0
2	A	10	0	0	3	0
2	B	5	0	0	0	0
2	C	10	0	0	2	0
2	D	5	0	0	0	0
2	E	5	0	0	2	0
2	G	5	0	0	1	0
2	H	5	0	0	0	0
2	I	5	0	0	0	0
2	K	5	0	0	2	0
2	L	5	0	0	0	0
All	All	18301	0	18506	1060	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 29.

All (1060) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:35:ARG:NH2	1:I:145:LEU:HD11	1.46	1.27
1:H:171:GLY:HA3	1:H:188:THR:O	1.60	1.01
1:F:23:PHE:HE1	1:F:128:ILE:HG12	1.28	0.97
1:I:34:GLN:HB3	1:I:139:VAL:HG21	1.43	0.97
1:A:82:GLN:HB2	1:D:196:MET:HE3	1.46	0.96
1:B:63:LYS:HE2	1:H:63:LYS:HE2	1.49	0.95
1:L:171:GLY:HA3	1:L:188:THR:O	1.69	0.93
1:I:35:ARG:HE	1:I:88:ILE:HD12	1.31	0.93
1:D:23:PHE:HE1	1:D:128:ILE:HG12	1.35	0.92
1:J:16:SER:HB2	1:J:109:THR:HG21	1.52	0.90
1:H:16:SER:HB2	1:H:109:THR:HG21	1.53	0.90
1:J:23:PHE:HE1	1:J:128:ILE:HG12	1.36	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:16:SER:HB2	1:F:109:THR:HG21	1.54	0.90
1:H:23:PHE:HE1	1:H:128:ILE:HG12	1.38	0.89
1:B:23:PHE:HE1	1:B:128:ILE:HG12	1.35	0.89
1:I:35:ARG:NH2	1:I:145:LEU:CD1	2.36	0.89
1:L:16:SER:HB2	1:L:109:THR:HG21	1.54	0.88
1:D:16:SER:HB2	1:D:109:THR:HG21	1.55	0.88
1:L:23:PHE:HE1	1:L:128:ILE:HG12	1.38	0.88
1:K:32:LEU:HD22	1:K:37:ILE:HG21	1.55	0.87
1:A:32:LEU:HD22	1:A:37:ILE:HG21	1.57	0.87
1:I:144:LEU:HD23	1:J:129:ARG:HH21	1.39	0.87
1:I:35:ARG:CZ	1:I:98:GLU:OE2	2.23	0.87
1:B:16:SER:HB2	1:B:109:THR:HG21	1.55	0.86
1:C:32:LEU:HD22	1:C:37:ILE:HG21	1.58	0.85
1:D:122:LYS:HD3	1:K:41:GLU:HG3	1.57	0.84
1:E:32:LEU:HD22	1:E:37:ILE:HG21	1.57	0.84
1:J:90:ASN:HB3	1:J:93:SER:HB3	1.60	0.84
1:G:32:LEU:HD22	1:G:37:ILE:HG21	1.58	0.83
1:L:90:ASN:HB3	1:L:93:SER:HB3	1.60	0.83
1:I:35:ARG:HH11	1:I:35:ARG:HG2	1.44	0.83
1:B:90:ASN:HB3	1:B:93:SER:HB3	1.60	0.83
1:I:35:ARG:HH22	1:I:145:LEU:HD11	1.43	0.83
1:D:173:GLN:HB2	1:D:187:THR:HG22	1.60	0.83
1:D:90:ASN:HB3	1:D:93:SER:HB3	1.59	0.83
1:B:200:LYS:HB3	1:B:202:PRO:HD3	1.61	0.83
1:I:138:THR:HA	1:I:141:PHE:CE1	2.14	0.82
1:I:144:LEU:CD2	1:J:129:ARG:HH21	1.93	0.82
1:B:196:MET:HE3	1:C:82:GLN:HB2	1.61	0.82
1:B:173:GLN:HB2	1:B:187:THR:HG22	1.60	0.81
1:L:48:LYS:HG2	1:L:49:TYR:CD1	2.15	0.81
1:F:90:ASN:HB3	1:F:93:SER:HB3	1.61	0.81
1:D:200:LYS:HB3	1:D:202:PRO:HD3	1.62	0.80
1:D:171:GLY:HA3	1:D:188:THR:O	1.81	0.80
1:I:35:ARG:HB3	1:I:35:ARG:NH1	1.96	0.80
1:F:23:PHE:CE1	1:F:128:ILE:HG12	2.16	0.80
1:J:173:GLN:HB2	1:J:187:THR:HG22	1.64	0.80
1:I:38:TYR:HB3	1:I:39:PRO:CD	2.12	0.80
1:J:200:LYS:HB3	1:J:202:PRO:HD3	1.64	0.80
1:F:173:GLN:HB2	1:F:187:THR:HG22	1.65	0.79
1:H:75:TRP:CH2	1:H:190:ILE:HD11	2.17	0.79
1:E:172:PRO:HB2	1:E:174:PHE:CE1	2.17	0.79
1:I:124:ILE:HD13	1:I:188:THR:HG22	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ILE:HD13	1:A:188:THR:HG22	1.61	0.79
1:I:35:ARG:CZ	1:I:98:GLU:CD	2.51	0.79
1:H:91:ILE:HD13	1:H:150:SER:HB3	1.64	0.79
1:J:91:ILE:HD13	1:J:150:SER:HB3	1.63	0.79
1:B:48:LYS:HG2	1:B:49:TYR:CD1	2.17	0.79
1:H:200:LYS:HB3	1:H:202:PRO:HD3	1.63	0.79
1:C:124:ILE:HD13	1:C:188:THR:HG22	1.64	0.79
1:I:133:ARG:HD3	1:J:34:GLN:HA	1.65	0.79
1:A:26:PHE:CE1	1:A:48:LYS:HG2	2.18	0.79
1:D:48:LYS:HG2	1:D:49:TYR:CD1	2.18	0.79
1:G:124:ILE:HD13	1:G:188:THR:HG22	1.65	0.79
1:E:65:LEU:O	1:E:69:VAL:HG23	1.83	0.78
1:B:171:GLY:HA3	1:B:188:THR:O	1.83	0.78
1:H:90:ASN:HB3	1:H:93:SER:HB3	1.63	0.78
1:L:91:ILE:HD13	1:L:150:SER:HB3	1.65	0.78
1:I:133:ARG:CD	1:J:34:GLN:HA	2.14	0.78
1:C:49:TYR:O	1:C:129:ARG:HD3	1.84	0.78
1:E:26:PHE:CE1	1:E:48:LYS:HG2	2.19	0.78
1:E:49:TYR:O	1:E:129:ARG:HD3	1.83	0.77
1:F:91:ILE:HD13	1:F:150:SER:HB3	1.67	0.77
1:K:124:ILE:HD13	1:K:188:THR:HG22	1.66	0.77
1:G:26:PHE:CE1	1:G:48:LYS:HG2	2.19	0.77
1:I:34:GLN:CB	1:I:139:VAL:HG21	2.13	0.77
1:B:46:VAL:HA	1:H:46:VAL:HA	1.66	0.77
1:G:37:ILE:HD11	1:G:88:ILE:HD13	1.67	0.76
1:C:10:SER:HA	1:C:117:ARG:O	1.85	0.76
1:L:75:TRP:CH2	1:L:190:ILE:HD11	2.19	0.76
1:K:172:PRO:HB2	1:K:174:PHE:CE1	2.21	0.76
1:G:170:SER:O	1:J:198:ALA:HB3	1.84	0.76
1:E:124:ILE:HD13	1:E:188:THR:HG22	1.67	0.76
1:K:26:PHE:CE1	1:K:48:LYS:HG2	2.21	0.76
1:H:48:LYS:HG2	1:H:49:TYR:CD1	2.20	0.76
1:A:56:THR:HG21	1:A:61:LEU:HD23	1.68	0.76
1:A:72:LEU:HD11	1:A:155:ILE:HD11	1.68	0.76
1:A:37:ILE:HD11	1:A:88:ILE:HD13	1.68	0.76
1:C:26:PHE:CE1	1:C:48:LYS:HG2	2.19	0.75
1:K:56:THR:HG21	1:K:61:LEU:HD23	1.66	0.75
1:E:56:THR:HG21	1:E:61:LEU:HD23	1.69	0.75
1:I:51:LEU:HD22	1:J:40:SER:OG	1.86	0.75
1:J:171:GLY:HA3	1:J:188:THR:O	1.86	0.75
1:B:75:TRP:CH2	1:B:190:ILE:HD11	2.21	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:48:LYS:HG2	1:J:49:TYR:CD1	2.20	0.75
1:J:75:TRP:CH2	1:J:190:ILE:HD11	2.20	0.75
1:I:10:SER:HA	1:I:117:ARG:O	1.87	0.75
1:A:10:SER:HA	1:A:117:ARG:O	1.87	0.75
1:K:49:TYR:O	1:K:129:ARG:HD3	1.87	0.75
1:I:49:TYR:O	1:I:129:ARG:HD3	1.86	0.74
1:F:171:GLY:HA3	1:F:188:THR:O	1.87	0.74
1:A:129:ARG:NH2	2:A:207:SO4:O3	2.21	0.74
1:A:49:TYR:O	1:A:129:ARG:HD3	1.87	0.74
1:D:75:TRP:CH2	1:D:190:ILE:HD11	2.23	0.74
1:C:41:GLU:HG3	1:L:122:LYS:HD3	1.68	0.74
1:I:35:ARG:NE	1:I:98:GLU:OE2	2.21	0.74
1:E:26:PHE:HE1	1:E:48:LYS:HG2	1.53	0.74
1:D:91:ILE:HD13	1:D:150:SER:HB3	1.69	0.74
1:L:62:ILE:O	1:L:66:ASN:HB2	1.88	0.73
1:D:33:TYR:HB2	1:D:43:PHE:CE2	2.23	0.73
1:L:75:TRP:CD1	1:L:161:LEU:HD21	2.22	0.73
1:G:49:TYR:O	1:G:129:ARG:HD3	1.85	0.73
1:G:56:THR:HG21	1:G:61:LEU:HD23	1.69	0.73
1:A:82:GLN:CB	1:D:196:MET:HE3	2.18	0.73
1:I:140:THR:HB	1:J:52:THR:O	1.89	0.73
1:G:26:PHE:HE1	1:G:48:LYS:HG2	1.54	0.73
1:F:75:TRP:CH2	1:F:190:ILE:HD11	2.23	0.73
1:G:65:LEU:O	1:G:69:VAL:HG23	1.89	0.73
1:C:37:ILE:HD11	1:C:88:ILE:HD13	1.71	0.73
1:G:72:LEU:HD11	1:G:155:ILE:HD11	1.71	0.73
1:C:72:LEU:HD11	1:C:155:ILE:HD11	1.69	0.73
1:I:136:THR:O	1:I:139:VAL:HG12	1.89	0.72
1:G:10:SER:HA	1:G:117:ARG:O	1.88	0.72
1:I:26:PHE:CE1	1:I:48:LYS:HG2	2.24	0.72
1:K:10:SER:HA	1:K:117:ARG:O	1.88	0.72
1:C:38:TYR:CE1	1:C:61:LEU:HD22	2.24	0.72
1:J:23:PHE:CE1	1:J:128:ILE:HG12	2.24	0.72
1:B:33:TYR:HB2	1:B:43:PHE:CE2	2.25	0.72
1:G:154:LEU:HD22	1:J:199:TYR:O	1.89	0.72
1:I:65:LEU:O	1:I:69:VAL:HG23	1.89	0.72
1:E:37:ILE:HD11	1:E:88:ILE:HD13	1.71	0.72
1:H:201:ILE:HG22	1:H:201:ILE:O	1.90	0.72
1:I:35:ARG:CG	1:I:35:ARG:HH11	2.02	0.72
1:D:201:ILE:O	1:D:201:ILE:HG22	1.89	0.71
1:I:35:ARG:HH21	1:I:145:LEU:HD11	1.52	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:62:ILE:O	1:J:66:ASN:HB2	1.90	0.71
1:K:65:LEU:O	1:K:69:VAL:HG23	1.91	0.71
1:H:198:ALA:HB3	1:I:170:SER:O	1.91	0.71
1:E:72:LEU:HD11	1:E:155:ILE:HD11	1.71	0.71
1:F:47:GLN:HB3	1:I:146:GLU:C	2.11	0.71
1:G:38:TYR:CE1	1:G:61:LEU:HD22	2.25	0.71
1:I:134:GLN:O	1:I:138:THR:HG23	1.91	0.71
1:A:26:PHE:HE1	1:A:48:LYS:HG2	1.55	0.71
1:H:201:ILE:HD11	1:I:167:TRP:CZ3	2.25	0.71
1:B:201:ILE:O	1:B:201:ILE:HG22	1.90	0.71
1:F:120:SER:HB2	1:I:58:ASP:CG	2.11	0.71
1:H:201:ILE:HG12	1:I:167:TRP:CZ3	2.25	0.71
1:F:47:GLN:HG3	1:F:47:GLN:O	1.91	0.71
1:I:72:LEU:HD11	1:I:155:ILE:HD11	1.73	0.71
1:C:26:PHE:HE1	1:C:48:LYS:HG2	1.54	0.71
1:I:35:ARG:HB3	1:I:35:ARG:CZ	2.21	0.70
1:E:38:TYR:CE1	1:E:61:LEU:HD22	2.26	0.70
1:K:72:LEU:HD11	1:K:155:ILE:HD11	1.73	0.70
1:A:59:LEU:HD21	1:E:55:VAL:HG21	1.72	0.70
1:F:62:ILE:O	1:F:66:ASN:HB2	1.91	0.70
1:J:201:ILE:O	1:J:201:ILE:HG22	1.90	0.70
1:C:58:ASP:O	1:C:62:ILE:HG13	1.92	0.70
1:B:63:LYS:CE	1:H:63:LYS:HE2	2.20	0.70
1:L:23:PHE:CE1	1:L:128:ILE:HG12	2.26	0.70
1:K:26:PHE:HE1	1:K:48:LYS:HG2	1.56	0.70
1:D:62:ILE:O	1:D:66:ASN:HB2	1.91	0.70
1:C:33:TYR:CE1	1:C:54:LEU:HD22	2.27	0.70
1:F:47:GLN:NE2	1:I:38:TYR:HE2	1.90	0.70
1:F:147:VAL:HG23	1:F:148:SER:O	1.92	0.70
1:D:42:THR:HB	1:D:43:PHE:CD1	2.27	0.69
1:J:147:VAL:HG23	1:J:148:SER:O	1.92	0.69
1:C:65:LEU:O	1:C:69:VAL:HG23	1.92	0.69
1:J:33:TYR:HB2	1:J:43:PHE:CE2	2.27	0.69
1:B:147:VAL:HG23	1:B:148:SER:O	1.91	0.69
1:B:62:ILE:O	1:B:66:ASN:HB2	1.92	0.69
1:D:23:PHE:CE1	1:D:128:ILE:HG12	2.23	0.69
1:I:144:LEU:CD2	1:J:129:ARG:NH2	2.55	0.69
1:G:170:SER:N	1:J:198:ALA:O	2.23	0.69
1:F:42:THR:HB	1:F:43:PHE:CD1	2.28	0.68
1:H:201:ILE:HG12	1:I:167:TRP:CE3	2.28	0.68
1:H:62:ILE:O	1:H:66:ASN:HB2	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:56:THR:HG21	1:I:61:LEU:HD23	1.73	0.68
1:I:34:GLN:HG2	1:J:45:ARG:HH22	1.57	0.68
1:K:37:ILE:HD11	1:K:88:ILE:HD13	1.73	0.68
1:D:147:VAL:HG23	1:D:148:SER:O	1.94	0.68
1:K:38:TYR:CE1	1:K:61:LEU:HD22	2.29	0.68
1:B:23:PHE:CE1	1:B:128:ILE:HG12	2.23	0.68
1:H:42:THR:HB	1:H:43:PHE:CD1	2.29	0.68
1:C:56:THR:HG21	1:C:61:LEU:HD23	1.74	0.68
1:H:147:VAL:HG23	1:H:148:SER:O	1.94	0.68
1:E:31:ILE:O	1:E:35:ARG:HG2	1.94	0.68
1:B:63:LYS:HE2	1:H:63:LYS:CE	2.23	0.67
1:L:48:LYS:HG2	1:L:49:TYR:CE1	2.28	0.67
1:B:48:LYS:HG2	1:B:49:TYR:CE1	2.29	0.67
1:I:34:GLN:OE1	1:I:136:THR:HG23	1.93	0.67
1:I:133:ARG:HD3	1:J:33:TYR:O	1.93	0.67
1:D:42:THR:HB	1:D:43:PHE:HD1	1.58	0.67
1:I:172:PRO:HB2	1:I:174:PHE:CE1	2.29	0.67
1:F:49:TYR:CE2	1:F:124:ILE:HG21	2.29	0.67
1:H:168:GLU:HB2	1:H:170:SER:O	1.95	0.67
1:H:74:ASP:O	1:H:77:TYR:HD2	1.77	0.67
1:D:63:LYS:HG3	1:L:59:LEU:HD21	1.74	0.67
1:K:83:LYS:HB2	1:K:156:TYR:HB2	1.76	0.67
1:L:74:ASP:O	1:L:77:TYR:HD2	1.78	0.67
1:E:141:PHE:HA	1:F:129:ARG:HG2	1.76	0.67
1:F:42:THR:HB	1:F:43:PHE:HD1	1.59	0.67
1:L:42:THR:HB	1:L:43:PHE:CD1	2.29	0.67
1:G:31:ILE:O	1:G:35:ARG:HG2	1.95	0.67
1:L:147:VAL:HG23	1:L:148:SER:O	1.94	0.67
1:H:33:TYR:HB2	1:H:43:PHE:CE2	2.29	0.66
1:B:91:ILE:HD13	1:B:150:SER:HB3	1.76	0.66
1:I:34:GLN:HG2	1:J:45:ARG:NH2	2.10	0.66
1:D:122:LYS:HA	1:K:41:GLU:HB2	1.76	0.66
1:A:129:ARG:NH1	2:A:207:SO4:O2	2.29	0.66
1:B:42:THR:HB	1:B:43:PHE:CD1	2.31	0.66
1:K:31:ILE:O	1:K:35:ARG:HG2	1.96	0.66
1:I:144:LEU:HD23	1:J:129:ARG:NH2	2.10	0.66
1:I:58:ASP:O	1:I:62:ILE:HG13	1.95	0.66
1:G:167:TRP:CZ3	1:J:201:ILE:HD11	2.30	0.66
1:J:130:SER:O	1:J:134:GLN:HG3	1.96	0.66
1:F:26:PHE:HE1	1:F:49:TYR:HE1	1.44	0.66
1:L:42:THR:HB	1:L:43:PHE:HD1	1.61	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:74:ASP:O	1:F:77:TYR:HD2	1.78	0.66
1:D:48:LYS:HG2	1:D:49:TYR:CE1	2.31	0.66
1:H:201:ILE:CD1	1:I:167:TRP:CZ3	2.78	0.66
1:H:146:GLU:HA	1:H:182:ARG:NH2	2.11	0.66
1:K:58:ASP:O	1:K:62:ILE:HG13	1.96	0.66
1:E:83:LYS:HB2	1:E:156:TYR:HB2	1.77	0.66
1:H:42:THR:HB	1:H:43:PHE:HD1	1.60	0.66
1:I:26:PHE:HE1	1:I:48:LYS:HG2	1.60	0.66
1:F:130:SER:O	1:F:134:GLN:HG3	1.95	0.66
1:B:74:ASP:O	1:B:77:TYR:HD2	1.79	0.66
1:E:172:PRO:CB	1:E:174:PHE:CE1	2.79	0.65
1:A:65:LEU:O	1:A:69:VAL:HG23	1.96	0.65
1:I:141:PHE:O	1:J:129:ARG:HD3	1.96	0.65
1:A:38:TYR:CE1	1:A:61:LEU:HD22	2.30	0.65
1:J:146:GLU:HA	1:J:182:ARG:NH2	2.12	0.65
1:G:172:PRO:HB2	1:G:174:PHE:CE1	2.31	0.65
1:C:172:PRO:HB2	1:C:174:PHE:CE1	2.31	0.65
1:J:42:THR:HB	1:J:43:PHE:CD1	2.32	0.65
1:L:146:GLU:HA	1:L:182:ARG:NH2	2.11	0.65
1:B:130:SER:O	1:B:134:GLN:HG3	1.97	0.65
1:B:146:GLU:HA	1:B:182:ARG:NH2	2.12	0.65
1:G:58:ASP:O	1:G:62:ILE:HG13	1.95	0.65
1:H:23:PHE:CE1	1:H:128:ILE:HG12	2.26	0.65
1:E:58:ASP:O	1:E:62:ILE:HG13	1.97	0.65
1:B:42:THR:HB	1:B:43:PHE:HD1	1.62	0.64
1:F:33:TYR:HB2	1:F:43:PHE:CE2	2.32	0.64
1:H:48:LYS:HG2	1:H:49:TYR:CE1	2.32	0.64
1:B:63:LYS:NZ	1:H:63:LYS:CE	2.60	0.64
1:L:33:TYR:HB2	1:L:43:PHE:CE2	2.32	0.64
1:K:33:TYR:CE1	1:K:54:LEU:HD22	2.32	0.64
1:B:63:LYS:CE	1:H:63:LYS:CE	2.75	0.64
1:H:130:SER:O	1:H:134:GLN:HG3	1.97	0.64
1:L:168:GLU:HB2	1:L:170:SER:O	1.96	0.64
1:D:74:ASP:O	1:D:77:TYR:HD2	1.79	0.64
1:D:130:SER:O	1:D:134:GLN:HG3	1.97	0.64
1:L:130:SER:O	1:L:134:GLN:HG3	1.98	0.64
1:A:83:LYS:HB2	1:A:156:TYR:HB2	1.80	0.64
1:D:46:VAL:HA	1:L:46:VAL:HA	1.80	0.64
1:H:163:VAL:HG22	1:H:190:ILE:HD12	1.79	0.64
1:I:83:LYS:HB2	1:I:156:TYR:HB2	1.79	0.64
1:A:172:PRO:HB2	1:A:174:PHE:CE1	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:36:GLY:O	1:I:37:ILE:HG22	1.97	0.64
1:F:146:GLU:HA	1:F:182:ARG:NH2	2.12	0.64
1:J:42:THR:HB	1:J:43:PHE:HD1	1.63	0.63
1:J:74:ASP:O	1:J:77:TYR:HD2	1.79	0.63
1:F:168:GLU:O	1:F:169:GLU:HB3	1.98	0.63
1:C:31:ILE:O	1:C:35:ARG:HG2	1.99	0.63
1:A:129:ARG:HH12	1:A:133:ARG:NH2	1.97	0.63
1:A:33:TYR:CE1	1:A:54:LEU:HD22	2.34	0.63
1:A:58:ASP:O	1:A:62:ILE:HG13	1.97	0.63
1:G:83:LYS:HB2	1:G:156:TYR:HB2	1.81	0.62
1:I:35:ARG:NH2	1:I:98:GLU:CD	2.52	0.62
1:J:168:GLU:O	1:J:169:GLU:HB3	1.98	0.62
1:E:33:TYR:CE1	1:E:54:LEU:HD22	2.34	0.62
1:I:34:GLN:HB3	1:I:139:VAL:CG2	2.27	0.62
1:C:129:ARG:NH2	2:C:206:SO4:O4	2.30	0.62
1:C:72:LEU:HD11	1:C:155:ILE:CD1	2.29	0.62
1:D:146:GLU:HA	1:D:182:ARG:NH2	2.14	0.62
1:I:33:TYR:CE1	1:I:54:LEU:HD22	2.33	0.62
1:I:51:LEU:HD21	1:I:129:ARG:HG2	1.82	0.62
1:J:48:LYS:HG2	1:J:49:TYR:CE1	2.34	0.62
1:K:30:SER:O	1:K:34:GLN:HG3	1.99	0.62
1:L:163:VAL:HG22	1:L:190:ILE:HD12	1.80	0.62
1:I:85:VAL:HB	1:I:154:LEU:HB2	1.82	0.62
1:A:72:LEU:HD11	1:A:155:ILE:CD1	2.30	0.61
1:B:63:LYS:CE	1:H:63:LYS:NZ	2.63	0.61
1:G:51:LEU:HD21	1:G:129:ARG:HG2	1.82	0.61
1:C:33:TYR:CZ	1:C:54:LEU:HD22	2.35	0.61
1:H:168:GLU:O	1:H:169:GLU:HB3	2.01	0.61
1:A:31:ILE:O	1:A:35:ARG:HG2	2.00	0.61
1:C:83:LYS:HB2	1:C:156:TYR:HB2	1.83	0.61
1:E:51:LEU:HD21	1:E:129:ARG:HG2	1.83	0.61
1:L:168:GLU:O	1:L:169:GLU:HB3	2.00	0.61
1:K:172:PRO:CB	1:K:174:PHE:CE1	2.83	0.60
1:E:33:TYR:CD2	1:E:54:LEU:HD13	2.35	0.60
1:I:35:ARG:NE	1:I:88:ILE:HD12	2.10	0.60
1:A:85:VAL:HB	1:A:154:LEU:HB2	1.84	0.60
1:F:46:VAL:HA	1:I:146:GLU:HB2	1.81	0.60
1:G:33:TYR:CE1	1:G:54:LEU:HD22	2.35	0.60
1:L:14:ARG:HA	1:L:105:GLU:HB3	1.84	0.60
1:G:129:ARG:HH12	1:G:133:ARG:NH2	2.00	0.60
1:C:85:VAL:HB	1:C:154:LEU:HB2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:133:ARG:HD3	1:H:33:TYR:O	2.02	0.60
1:L:14:ARG:HB3	1:L:105:GLU:HG2	1.83	0.59
1:I:35:ARG:CB	1:I:35:ARG:NH1	2.64	0.59
1:G:129:ARG:NH1	1:G:133:ARG:NH2	2.51	0.59
1:A:100:TRP:CD2	1:A:197:VAL:HG23	2.37	0.59
1:C:129:ARG:HH12	1:C:133:ARG:NH2	2.00	0.59
1:K:51:LEU:HD21	1:K:129:ARG:HG2	1.85	0.59
1:A:30:SER:O	1:A:34:GLN:HG3	2.03	0.59
1:I:133:ARG:HD3	1:J:34:GLN:CA	2.33	0.58
1:H:201:ILE:CG1	1:I:167:TRP:CZ3	2.85	0.58
1:I:72:LEU:HD11	1:I:155:ILE:CD1	2.33	0.58
1:H:64:TYR:HE1	1:H:172:PRO:O	1.87	0.58
1:A:129:ARG:NH1	1:A:133:ARG:NH2	2.51	0.58
1:B:43:PHE:N	1:B:43:PHE:HD1	2.01	0.58
1:A:59:LEU:HD22	1:E:44:THR:HG22	1.85	0.58
1:C:33:TYR:CD2	1:C:54:LEU:HD13	2.38	0.58
1:B:159:LYS:O	1:B:159:LYS:HG3	2.02	0.58
1:G:72:LEU:HD11	1:G:155:ILE:CD1	2.33	0.58
1:A:59:LEU:HD22	1:E:44:THR:CG2	2.33	0.58
1:G:42:THR:HB	1:G:43:PHE:CD1	2.39	0.58
1:D:159:LYS:HG3	1:D:159:LYS:O	2.02	0.58
1:K:33:TYR:CZ	1:K:54:LEU:HD22	2.38	0.58
1:G:33:TYR:CD2	1:G:54:LEU:HD13	2.39	0.58
1:I:38:TYR:HB3	1:I:39:PRO:HD2	1.85	0.58
1:D:168:GLU:O	1:D:169:GLU:HB3	2.03	0.58
1:G:30:SER:O	1:G:34:GLN:HG3	2.04	0.58
1:G:91:ILE:HD13	1:G:150:SER:OG	2.04	0.58
1:B:200:LYS:HE3	1:C:170:SER:OG	2.03	0.58
1:B:196:MET:HE3	1:C:82:GLN:CB	2.33	0.58
1:E:72:LEU:HD11	1:E:155:ILE:CD1	2.33	0.58
1:C:91:ILE:HD13	1:C:150:SER:OG	2.04	0.57
1:L:162:VAL:HG23	1:L:162:VAL:O	2.04	0.57
1:I:91:ILE:HD13	1:I:150:SER:OG	2.03	0.57
1:E:133:ARG:HD3	1:F:33:TYR:O	2.04	0.57
1:E:91:ILE:HD13	1:E:150:SER:OG	2.04	0.57
1:A:91:ILE:HD13	1:A:150:SER:OG	2.03	0.57
1:F:14:ARG:HA	1:F:105:GLU:HB3	1.85	0.57
1:K:72:LEU:HD11	1:K:155:ILE:CD1	2.34	0.57
1:B:168:GLU:O	1:B:169:GLU:HB3	2.03	0.57
1:A:33:TYR:CD2	1:A:54:LEU:HD13	2.39	0.57
1:B:14:ARG:HA	1:B:105:GLU:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:159:LYS:O	1:F:159:LYS:HG3	2.04	0.57
1:K:91:ILE:HD13	1:K:150:SER:OG	2.05	0.57
1:F:49:TYR:CE1	1:F:128:ILE:HD11	2.40	0.56
1:E:172:PRO:HB3	1:E:174:PHE:CZ	2.40	0.56
1:C:30:SER:O	1:C:34:GLN:HG3	2.05	0.56
1:J:159:LYS:O	1:J:159:LYS:HG3	2.03	0.56
1:K:33:TYR:CD2	1:K:54:LEU:HD13	2.39	0.56
1:H:31:ILE:HD13	1:H:100:TRP:CD2	2.40	0.56
1:B:63:LYS:HZ3	1:H:63:LYS:CE	2.19	0.56
1:F:16:SER:O	1:F:19:ILE:HG22	2.04	0.56
1:I:144:LEU:HD21	1:J:129:ARG:NH2	2.20	0.56
1:D:91:ILE:CD1	1:D:150:SER:HB3	2.34	0.56
1:B:165:GLU:CD	1:B:165:GLU:H	2.09	0.56
1:H:14:ARG:HA	1:H:105:GLU:HB3	1.88	0.56
1:F:88:ILE:HG12	1:F:151:PHE:CE1	2.40	0.56
1:K:42:THR:HB	1:K:43:PHE:CD1	2.40	0.56
1:F:165:GLU:H	1:F:165:GLU:CD	2.07	0.56
1:H:16:SER:O	1:H:19:ILE:HG22	2.05	0.56
1:F:47:GLN:HB3	1:I:146:GLU:O	2.05	0.56
1:L:43:PHE:N	1:L:43:PHE:HD1	2.04	0.56
1:K:129:ARG:HH12	1:K:133:ARG:NH2	2.03	0.56
1:A:42:THR:HB	1:A:43:PHE:CD1	2.40	0.56
1:I:129:ARG:NH1	1:I:133:ARG:NH2	2.54	0.56
1:B:43:PHE:N	1:B:43:PHE:CD1	2.72	0.56
1:C:129:ARG:NH1	1:C:133:ARG:NH2	2.54	0.56
1:D:14:ARG:HA	1:D:105:GLU:HB3	1.87	0.56
1:H:8:HIS:CG	1:H:9:GLY:H	2.24	0.56
1:I:35:ARG:CG	1:I:35:ARG:NH1	2.64	0.56
1:D:179:GLU:CD	1:D:180:GLU:H	2.09	0.56
1:K:129:ARG:NH1	1:K:133:ARG:NH2	2.53	0.56
1:H:43:PHE:N	1:H:43:PHE:HD1	2.04	0.56
1:D:43:PHE:N	1:D:43:PHE:HD1	2.04	0.56
1:E:129:ARG:HH12	1:E:133:ARG:NH2	2.04	0.56
1:I:34:GLN:OE1	1:J:45:ARG:NH2	2.39	0.55
1:G:85:VAL:HB	1:G:154:LEU:HB2	1.88	0.55
1:I:31:ILE:HD13	1:I:100:TRP:CD2	2.41	0.55
1:I:139:VAL:HG13	1:I:140:THR:N	2.20	0.55
1:F:19:ILE:CG2	1:F:20:VAL:N	2.70	0.55
1:A:133:ARG:HD3	1:B:33:TYR:O	2.05	0.55
1:A:33:TYR:CZ	1:A:54:LEU:HD22	2.41	0.55
1:E:33:TYR:CZ	1:E:54:LEU:HD22	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:90:ASN:CG	1:D:93:SER:HB2	2.27	0.55
1:A:51:LEU:HD21	1:A:129:ARG:HG2	1.87	0.55
1:I:172:PRO:CB	1:I:174:PHE:CE1	2.89	0.55
1:A:91:ILE:HG22	1:A:148:SER:O	2.06	0.55
1:C:183:LEU:HB2	1:C:195:SER:O	2.07	0.55
1:F:136:THR:O	1:F:139:VAL:HG13	2.07	0.55
1:C:31:ILE:HD13	1:C:100:TRP:CD2	2.42	0.55
1:H:173:GLN:HB2	1:H:187:THR:HG22	1.89	0.55
1:E:129:ARG:NH1	1:E:133:ARG:NH2	2.55	0.55
1:B:31:ILE:HD13	1:B:100:TRP:CD2	2.42	0.55
1:J:179:GLU:CD	1:J:180:GLU:H	2.10	0.55
1:I:40:SER:O	1:I:43:PHE:O	2.25	0.55
1:K:164:PRO:O	1:K:167:TRP:HD1	1.89	0.55
1:I:24:PHE:CE2	1:I:84:LEU:HG	2.43	0.55
1:K:85:VAL:HB	1:K:154:LEU:HB2	1.89	0.55
1:A:59:LEU:CD2	1:E:55:VAL:CG2	2.85	0.54
1:L:31:ILE:HD13	1:L:100:TRP:CD2	2.42	0.54
1:D:43:PHE:N	1:D:43:PHE:CD1	2.75	0.54
1:A:33:TYR:CE2	1:A:54:LEU:HD13	2.42	0.54
1:L:179:GLU:CD	1:L:180:GLU:H	2.11	0.54
1:C:108:LYS:HD2	1:C:111:LYS:NZ	2.22	0.54
1:A:164:PRO:O	1:A:167:TRP:HD1	1.91	0.54
1:F:31:ILE:HD13	1:F:100:TRP:CD2	2.43	0.54
1:E:164:PRO:O	1:E:167:TRP:HD1	1.91	0.54
1:E:33:TYR:CE2	1:E:54:LEU:HD13	2.43	0.54
1:A:108:LYS:HD2	1:A:111:LYS:NZ	2.22	0.54
1:J:31:ILE:HD13	1:J:100:TRP:CD2	2.43	0.54
1:K:31:ILE:HD13	1:K:100:TRP:CD2	2.43	0.54
1:E:85:VAL:HB	1:E:154:LEU:HB2	1.89	0.54
1:F:19:ILE:HG23	1:F:20:VAL:N	2.23	0.54
1:L:43:PHE:N	1:L:43:PHE:CD1	2.74	0.54
1:G:164:PRO:O	1:G:167:TRP:HD1	1.91	0.54
1:E:42:THR:HB	1:E:43:PHE:CD1	2.42	0.54
1:I:129:ARG:HH12	1:I:133:ARG:NH2	2.06	0.54
1:E:31:ILE:HD13	1:E:100:TRP:CD2	2.43	0.54
1:K:100:TRP:CD2	1:K:197:VAL:HG23	2.42	0.54
1:H:19:ILE:CG2	1:H:20:VAL:N	2.71	0.54
1:C:41:GLU:HB2	1:L:122:LYS:HA	1.89	0.54
1:C:164:PRO:O	1:C:167:TRP:HD1	1.90	0.54
1:G:167:TRP:HZ3	1:J:201:ILE:HD11	1.72	0.54
1:H:43:PHE:N	1:H:43:PHE:CD1	2.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:SER:O	1:E:34:GLN:HG3	2.08	0.53
1:I:164:PRO:O	1:I:167:TRP:HD1	1.92	0.53
1:A:59:LEU:HD21	1:E:55:VAL:CG2	2.39	0.53
1:G:33:TYR:CZ	1:G:54:LEU:HD22	2.43	0.53
1:F:179:GLU:CD	1:F:180:GLU:H	2.11	0.53
1:B:179:GLU:CD	1:B:180:GLU:H	2.11	0.53
1:C:28:ILE:HD13	1:C:153:LEU:HD21	1.90	0.53
1:L:16:SER:O	1:L:19:ILE:HG22	2.09	0.53
1:L:90:ASN:CB	1:L:93:SER:HB3	2.37	0.53
1:A:183:LEU:HB2	1:A:195:SER:O	2.09	0.53
1:F:91:ILE:CD1	1:F:150:SER:HB3	2.39	0.53
1:G:51:LEU:HD21	1:G:129:ARG:CG	2.38	0.53
1:K:33:TYR:CE2	1:K:54:LEU:HD13	2.44	0.53
1:D:165:GLU:CD	1:D:165:GLU:H	2.11	0.53
1:C:42:THR:HB	1:C:43:PHE:CD1	2.43	0.53
1:J:90:ASN:CB	1:J:93:SER:HB3	2.37	0.53
1:I:33:TYR:CZ	1:I:54:LEU:HD22	2.43	0.53
1:F:120:SER:CB	1:I:58:ASP:OD1	2.57	0.53
1:D:63:LYS:HE2	1:L:63:LYS:HE2	1.91	0.52
1:F:47:GLN:HE21	1:I:38:TYR:HE2	1.56	0.52
1:I:100:TRP:CD2	1:I:197:VAL:HG23	2.44	0.52
1:J:14:ARG:HA	1:J:105:GLU:HB3	1.91	0.52
1:L:19:ILE:CG2	1:L:20:VAL:N	2.72	0.52
1:G:31:ILE:HD13	1:G:100:TRP:CD2	2.44	0.52
1:G:64:TYR:C	1:G:64:TYR:CD1	2.83	0.52
1:I:34:GLN:CG	1:J:45:ARG:HH22	2.22	0.52
1:B:90:ASN:CG	1:B:93:SER:HB2	2.30	0.52
1:C:100:TRP:CD2	1:C:197:VAL:HG23	2.44	0.52
1:K:91:ILE:HG22	1:K:148:SER:O	2.10	0.52
1:L:175:ILE:HG22	1:L:176:THR:N	2.24	0.52
1:L:168:GLU:C	1:L:170:SER:H	2.13	0.52
1:G:54:LEU:HD12	1:G:54:LEU:N	2.24	0.52
1:A:43:PHE:N	1:A:43:PHE:CD1	2.77	0.52
1:J:43:PHE:CD1	1:J:43:PHE:N	2.77	0.52
1:G:100:TRP:CD2	1:G:197:VAL:HG23	2.44	0.52
1:A:172:PRO:CB	1:A:174:PHE:CE1	2.93	0.52
1:C:129:ARG:NH1	2:C:206:SO4:O1	2.40	0.52
1:F:43:PHE:CD1	1:F:43:PHE:N	2.77	0.52
1:C:43:PHE:N	1:C:43:PHE:CD1	2.78	0.52
1:K:64:TYR:C	1:K:64:TYR:CD1	2.83	0.52
1:J:43:PHE:HD1	1:J:43:PHE:N	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:49:TYR:HD2	1:G:125:GLN:HG3	1.75	0.52
1:C:33:TYR:CE2	1:C:54:LEU:HD13	2.44	0.52
1:I:24:PHE:CD2	1:I:84:LEU:HG	2.44	0.52
1:I:33:TYR:CD2	1:I:54:LEU:HD13	2.45	0.52
1:I:88:ILE:HB	1:I:98:GLU:HB3	1.92	0.52
1:K:88:ILE:HB	1:K:98:GLU:HB3	1.90	0.52
1:I:49:TYR:HD2	1:I:125:GLN:HG3	1.75	0.51
1:E:141:PHE:HA	1:F:129:ARG:CG	2.40	0.51
1:C:172:PRO:CB	1:C:174:PHE:CE1	2.93	0.51
1:G:91:ILE:HG22	1:G:148:SER:O	2.10	0.51
1:H:174:PHE:CE1	1:H:186:PHE:HB3	2.44	0.51
1:L:88:ILE:HG12	1:L:151:PHE:CE1	2.46	0.51
1:J:19:ILE:CG2	1:J:20:VAL:N	2.73	0.51
1:G:172:PRO:CB	1:G:174:PHE:CE1	2.92	0.51
1:C:88:ILE:HB	1:C:98:GLU:HB3	1.92	0.51
1:E:49:TYR:HD2	1:E:125:GLN:HG3	1.75	0.51
1:G:24:PHE:O	1:G:28:ILE:HG13	2.10	0.51
1:J:91:ILE:CD1	1:J:150:SER:HB3	2.36	0.51
1:C:133:ARG:HD3	1:D:33:TYR:O	2.09	0.51
1:A:24:PHE:O	1:A:28:ILE:HG13	2.10	0.51
1:L:197:VAL:HG11	1:L:199:TYR:OH	2.08	0.51
1:B:122:LYS:HD3	1:G:41:GLU:HG3	1.93	0.51
1:C:49:TYR:HD2	1:C:125:GLN:HG3	1.75	0.51
1:F:43:PHE:HD1	1:F:43:PHE:N	2.07	0.51
1:G:43:PHE:N	1:G:43:PHE:CD1	2.78	0.51
1:G:88:ILE:HB	1:G:98:GLU:HB3	1.93	0.51
1:H:179:GLU:CD	1:H:180:GLU:H	2.14	0.51
1:J:163:VAL:HG13	1:J:163:VAL:O	2.11	0.51
1:E:64:TYR:C	1:E:64:TYR:CD1	2.83	0.51
1:A:49:TYR:HD2	1:A:125:GLN:HG3	1.75	0.51
1:H:168:GLU:C	1:H:170:SER:H	2.13	0.51
1:E:141:PHE:O	1:F:129:ARG:HD3	2.10	0.51
1:H:8:HIS:CG	1:H:9:GLY:N	2.79	0.51
1:B:163:VAL:HG13	1:B:163:VAL:O	2.11	0.51
1:I:64:TYR:CD1	1:I:64:TYR:C	2.84	0.51
1:I:139:VAL:HA	1:I:142:LEU:HD12	1.93	0.51
1:D:90:ASN:ND2	1:D:93:SER:HB2	2.25	0.51
1:I:51:LEU:CD2	1:J:40:SER:OG	2.58	0.51
1:A:31:ILE:HD13	1:A:100:TRP:CD2	2.45	0.51
1:D:32:LEU:HD13	1:D:61:LEU:HD21	1.93	0.51
1:A:88:ILE:HB	1:A:98:GLU:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:172:PRO:HB3	1:K:174:PHE:CZ	2.46	0.51
1:K:49:TYR:HD2	1:K:125:GLN:HG3	1.76	0.51
1:E:100:TRP:CD2	1:E:197:VAL:HG23	2.46	0.51
1:C:78:LYS:O	1:C:79:CYS:HB2	2.09	0.51
1:A:64:TYR:C	1:A:64:TYR:CD1	2.85	0.51
1:E:88:ILE:HB	1:E:98:GLU:HB3	1.93	0.50
1:K:51:LEU:HD21	1:K:129:ARG:CG	2.41	0.50
1:C:24:PHE:CE2	1:C:84:LEU:HG	2.46	0.50
1:C:91:ILE:HG22	1:C:148:SER:O	2.11	0.50
1:B:88:ILE:HG12	1:B:151:PHE:CE1	2.46	0.50
1:E:178:SER:O	1:E:201:ILE:HD11	2.11	0.50
1:L:75:TRP:CG	1:L:161:LEU:HD21	2.47	0.50
1:C:24:PHE:O	1:C:28:ILE:HG13	2.12	0.50
1:E:24:PHE:O	1:E:28:ILE:HG13	2.10	0.50
1:L:135:ILE:O	1:L:139:VAL:HG12	2.11	0.50
1:H:136:THR:O	1:H:139:VAL:HG13	2.12	0.50
1:I:35:ARG:NH2	1:I:98:GLU:OE1	2.44	0.50
1:L:19:ILE:HG23	1:L:20:VAL:N	2.26	0.50
1:G:170:SER:O	1:J:198:ALA:CB	2.58	0.50
1:K:78:LYS:O	1:K:79:CYS:HB2	2.11	0.50
1:B:63:LYS:CE	1:H:63:LYS:HZ3	2.23	0.50
1:E:51:LEU:HD21	1:E:129:ARG:CG	2.42	0.50
1:K:43:PHE:N	1:K:43:PHE:CD1	2.80	0.50
1:E:43:PHE:CD1	1:E:43:PHE:N	2.77	0.50
1:L:197:VAL:CG1	1:L:199:TYR:CZ	2.94	0.50
1:H:91:ILE:CD1	1:H:150:SER:HB3	2.37	0.50
1:L:75:TRP:CD1	1:L:161:LEU:HD11	2.46	0.50
1:B:155:ILE:HB	1:B:190:ILE:HG12	1.93	0.50
1:D:137:ALA:O	1:D:140:THR:HB	2.11	0.50
1:D:90:ASN:CB	1:D:93:SER:HB3	2.35	0.50
1:F:174:PHE:CE1	1:F:186:PHE:HB3	2.47	0.50
1:I:164:PRO:HB2	1:I:167:TRP:HE1	1.76	0.50
1:C:27:GLY:O	1:C:31:ILE:HG13	2.12	0.50
1:K:164:PRO:HB2	1:K:167:TRP:HE1	1.77	0.50
1:K:32:LEU:HD22	1:K:37:ILE:CG2	2.36	0.50
1:H:173:GLN:CB	1:H:187:THR:HG22	2.41	0.50
1:G:78:LYS:O	1:G:79:CYS:HB2	2.12	0.50
1:F:16:SER:O	1:F:20:VAL:HG23	2.12	0.50
1:G:33:TYR:CE2	1:G:54:LEU:HD13	2.47	0.50
1:D:47:GLN:HB2	1:D:51:LEU:O	2.12	0.50
1:K:178:SER:O	1:K:201:ILE:HD11	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:34:GLN:HB3	1:I:139:VAL:HG11	1.93	0.49
1:D:19:ILE:CG2	1:D:20:VAL:N	2.75	0.49
1:E:129:ARG:O	1:E:133:ARG:HB2	2.12	0.49
1:B:201:ILE:HD11	1:C:167:TRP:CZ3	2.47	0.49
1:E:164:PRO:HB2	1:E:167:TRP:HE1	1.76	0.49
1:H:88:ILE:HG12	1:H:151:PHE:CE1	2.47	0.49
1:I:51:LEU:HD21	1:I:129:ARG:CG	2.41	0.49
1:I:91:ILE:HG22	1:I:148:SER:O	2.12	0.49
1:E:24:PHE:CE2	1:E:84:LEU:HG	2.48	0.49
1:J:88:ILE:HG12	1:J:151:PHE:CE1	2.46	0.49
1:J:174:PHE:CE1	1:J:186:PHE:HB3	2.46	0.49
1:I:40:SER:O	1:I:43:PHE:N	2.45	0.49
1:C:51:LEU:HD21	1:C:129:ARG:HG2	1.93	0.49
1:G:183:LEU:HB2	1:G:195:SER:O	2.12	0.49
1:A:51:LEU:HD21	1:A:129:ARG:CG	2.42	0.49
1:I:24:PHE:O	1:I:28:ILE:HG13	2.13	0.49
1:A:24:PHE:CE2	1:A:84:LEU:HG	2.48	0.49
1:B:174:PHE:CE1	1:B:186:PHE:HB3	2.48	0.49
1:H:19:ILE:HG23	1:H:20:VAL:N	2.27	0.49
1:C:32:LEU:HD22	1:C:37:ILE:CG2	2.38	0.49
1:B:16:SER:O	1:B:20:VAL:HG23	2.13	0.49
1:D:174:PHE:CE1	1:D:186:PHE:HB3	2.47	0.49
1:J:19:ILE:HG23	1:J:20:VAL:N	2.27	0.49
1:D:16:SER:O	1:D:20:VAL:HG23	2.13	0.49
1:F:33:TYR:CD2	1:F:54:LEU:HD23	2.47	0.49
1:C:24:PHE:CD2	1:C:84:LEU:HG	2.47	0.49
1:C:64:TYR:C	1:C:64:TYR:CD1	2.86	0.49
1:E:91:ILE:HG22	1:E:148:SER:O	2.12	0.49
1:A:178:SER:O	1:A:201:ILE:HD11	2.13	0.49
1:K:131:VAL:HG12	1:K:132:ILE:N	2.28	0.49
1:B:63:LYS:HZ3	1:H:63:LYS:HE3	1.78	0.49
1:B:19:ILE:CG2	1:B:20:VAL:N	2.75	0.49
1:D:122:LYS:CA	1:K:41:GLU:HB2	2.43	0.49
1:A:170:SER:OG	1:D:200:LYS:HE3	2.13	0.48
1:E:56:THR:CG2	1:E:61:LEU:HD23	2.42	0.48
1:G:24:PHE:CE2	1:G:84:LEU:HG	2.48	0.48
1:H:201:ILE:HD11	1:I:167:TRP:HZ3	1.74	0.48
1:H:32:LEU:HD13	1:H:61:LEU:HD21	1.95	0.48
1:I:142:LEU:O	1:I:143:PRO:C	2.51	0.48
1:J:136:THR:O	1:J:139:VAL:HG13	2.13	0.48
1:B:90:ASN:CB	1:B:93:SER:HB3	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:155:ILE:HB	1:F:190:ILE:HG12	1.94	0.48
1:C:38:TYR:CD1	1:C:61:LEU:HD22	2.48	0.48
1:E:183:LEU:HB2	1:E:195:SER:O	2.12	0.48
1:F:49:TYR:CD2	1:F:124:ILE:HG21	2.48	0.48
1:E:129:ARG:NH1	2:E:206:SO4:O4	2.40	0.48
1:D:31:ILE:HD13	1:D:100:TRP:CD2	2.47	0.48
1:K:56:THR:HG21	1:K:61:LEU:CD2	2.41	0.48
1:I:178:SER:O	1:I:201:ILE:HD11	2.13	0.48
1:H:167:TRP:CH2	1:H:172:PRO:HD2	2.49	0.48
1:F:47:GLN:NE2	1:I:38:TYR:CE2	2.76	0.48
1:G:152:ASP:OD2	1:J:200:LYS:HD3	2.14	0.48
1:A:164:PRO:HB2	1:A:167:TRP:HE1	1.78	0.48
1:C:43:PHE:O	1:L:121:GLN:HB3	2.14	0.48
1:I:35:ARG:CZ	1:I:98:GLU:OE1	2.62	0.48
1:H:167:TRP:CZ3	1:H:171:GLY:HA2	2.48	0.48
1:L:155:ILE:HB	1:L:190:ILE:HG12	1.96	0.48
1:G:164:PRO:HB2	1:G:167:TRP:HE1	1.78	0.48
1:A:24:PHE:CD2	1:A:84:LEU:HG	2.49	0.48
1:L:14:ARG:HB3	1:L:105:GLU:CB	2.44	0.48
1:F:50:GLY:O	1:I:148:SER:HB3	2.14	0.48
1:F:14:ARG:HB3	1:F:105:GLU:HG2	1.94	0.48
1:K:28:ILE:HD13	1:K:153:LEU:HD21	1.95	0.48
1:H:135:ILE:O	1:H:139:VAL:HG12	2.12	0.48
1:J:16:SER:O	1:J:19:ILE:HG22	2.14	0.48
1:C:54:LEU:HD12	1:C:54:LEU:N	2.29	0.48
1:F:135:ILE:O	1:F:139:VAL:HG12	2.13	0.48
1:F:163:VAL:O	1:F:163:VAL:HG13	2.14	0.48
1:F:14:ARG:HB3	1:F:105:GLU:CB	2.44	0.48
1:I:138:THR:HA	1:I:141:PHE:HE1	1.75	0.48
1:I:129:ARG:O	1:I:133:ARG:HB2	2.13	0.48
1:D:33:TYR:CD2	1:D:54:LEU:HD23	2.49	0.48
1:G:129:ARG:O	1:G:133:ARG:HB2	2.13	0.48
1:E:78:LYS:O	1:E:79:CYS:HB2	2.13	0.48
1:G:178:SER:O	1:G:201:ILE:HD11	2.14	0.48
1:E:172:PRO:CB	1:E:174:PHE:CZ	2.97	0.48
1:L:91:ILE:CD1	1:L:150:SER:HB3	2.39	0.48
1:H:14:ARG:HB3	1:H:105:GLU:CB	2.44	0.48
1:C:164:PRO:HB2	1:C:167:TRP:HE1	1.79	0.47
1:J:179:GLU:CG	1:J:180:GLU:N	2.77	0.47
1:I:138:THR:HG22	1:J:136:THR:HG21	1.96	0.47
1:J:33:TYR:CD2	1:J:54:LEU:HD23	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:56:THR:HG21	1:E:61:LEU:CD2	2.42	0.47
1:G:38:TYR:CD1	1:G:61:LEU:HD22	2.49	0.47
1:D:14:ARG:HB3	1:D:105:GLU:CB	2.44	0.47
1:D:179:GLU:CG	1:D:180:GLU:N	2.77	0.47
1:L:132:ILE:HD13	1:L:132:ILE:N	2.29	0.47
1:D:16:SER:O	1:D:19:ILE:HG22	2.14	0.47
1:J:135:ILE:O	1:J:139:VAL:HG12	2.14	0.47
1:I:38:TYR:HB3	1:I:39:PRO:HD3	1.95	0.47
1:C:51:LEU:HD21	1:C:129:ARG:CG	2.43	0.47
1:G:42:THR:HB	1:G:43:PHE:CE1	2.49	0.47
1:B:14:ARG:HB3	1:B:105:GLU:CB	2.44	0.47
1:L:33:TYR:CD2	1:L:54:LEU:HD23	2.50	0.47
1:J:155:ILE:HB	1:J:190:ILE:HG12	1.96	0.47
1:G:129:ARG:NH2	2:G:206:SO4:O1	2.47	0.47
1:E:54:LEU:HD12	1:E:54:LEU:N	2.30	0.47
1:K:84:LEU:HD13	1:K:85:VAL:N	2.29	0.47
1:F:48:LYS:HG3	1:F:49:TYR:CE1	2.49	0.47
1:L:167:TRP:CZ3	1:L:171:GLY:HA2	2.49	0.47
1:G:24:PHE:CD2	1:G:84:LEU:HG	2.49	0.47
1:L:179:GLU:CG	1:L:180:GLU:N	2.78	0.47
1:E:24:PHE:CD2	1:E:84:LEU:HG	2.49	0.47
1:B:135:ILE:O	1:B:139:VAL:HG12	2.15	0.47
1:L:24:PHE:O	1:L:28:ILE:HG22	2.15	0.47
1:D:163:VAL:O	1:D:163:VAL:HG13	2.15	0.47
1:A:129:ARG:O	1:A:133:ARG:HB2	2.14	0.47
1:E:49:TYR:O	1:E:129:ARG:CD	2.59	0.47
1:E:42:THR:HB	1:E:43:PHE:CE1	2.50	0.47
1:K:133:ARG:HD3	1:L:33:TYR:O	2.15	0.47
1:E:38:TYR:CD1	1:E:61:LEU:HD22	2.50	0.47
1:A:153:LEU:O	1:A:154:LEU:HD23	2.14	0.47
1:A:175:ILE:HB	1:A:178:SER:HB3	1.96	0.47
1:A:78:LYS:O	1:A:79:CYS:HB2	2.15	0.47
1:D:88:ILE:HG12	1:D:151:PHE:CE1	2.50	0.47
1:H:155:ILE:HB	1:H:190:ILE:HG12	1.96	0.47
1:A:42:THR:HB	1:A:43:PHE:CE1	2.50	0.47
1:L:137:ALA:O	1:L:140:THR:HB	2.15	0.47
1:A:56:THR:HG21	1:A:61:LEU:CD2	2.43	0.47
1:A:154:LEU:HD22	1:D:199:TYR:O	2.15	0.47
1:C:45:ARG:HD2	1:L:121:GLN:OE1	2.14	0.47
1:I:35:ARG:HH22	1:I:145:LEU:CD1	2.16	0.47
1:F:26:PHE:CE1	1:F:49:TYR:HE1	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:80:SER:HB2	1:H:161:LEU:HD22	1.97	0.47
1:D:155:ILE:HB	1:D:190:ILE:HG12	1.97	0.47
1:E:20:VAL:O	1:E:24:PHE:HD1	1.98	0.47
1:L:32:LEU:HD13	1:L:61:LEU:HD21	1.97	0.47
1:I:78:LYS:O	1:I:79:CYS:HB2	2.13	0.47
1:J:140:THR:HG22	1:J:141:PHE:CD1	2.50	0.47
1:H:39:PRO:HD3	1:H:181:VAL:HG22	1.98	0.46
1:K:42:THR:HB	1:K:43:PHE:CE1	2.50	0.46
1:D:180:GLU:OE1	1:D:180:GLU:HA	2.15	0.46
1:I:27:GLY:O	1:I:31:ILE:HG13	2.15	0.46
1:H:137:ALA:O	1:H:140:THR:HB	2.15	0.46
1:C:203:VAL:O	1:C:204:ASN:HB2	2.14	0.46
1:D:93:SER:OG	1:D:95:GLU:HG2	2.14	0.46
1:D:33:TYR:HB2	1:D:43:PHE:CD2	2.51	0.46
1:G:167:TRP:CZ3	1:J:201:ILE:CD1	2.98	0.46
1:K:175:ILE:HB	1:K:178:SER:HB3	1.97	0.46
1:C:11:ILE:HA	1:C:11:ILE:HD13	1.75	0.46
1:J:16:SER:O	1:J:20:VAL:HG23	2.15	0.46
1:K:24:PHE:CE2	1:K:84:LEU:HG	2.50	0.46
1:I:153:LEU:O	1:I:154:LEU:HD23	2.16	0.46
1:F:179:GLU:CG	1:F:180:GLU:N	2.78	0.46
1:E:175:ILE:HB	1:E:178:SER:HB3	1.98	0.46
1:G:175:ILE:HB	1:G:178:SER:HB3	1.98	0.46
1:C:131:VAL:HG12	1:C:132:ILE:N	2.29	0.46
1:K:183:LEU:HB2	1:K:195:SER:O	2.16	0.46
1:B:19:ILE:HG23	1:B:20:VAL:N	2.31	0.46
1:G:28:ILE:HD13	1:G:153:LEU:HD21	1.96	0.46
1:H:167:TRP:O	1:H:167:TRP:CE3	2.68	0.46
1:H:80:SER:CB	1:H:161:LEU:HD22	2.46	0.46
1:I:133:ARG:NH1	1:J:33:TYR:O	2.47	0.46
1:H:33:TYR:CD2	1:H:54:LEU:HD23	2.50	0.46
1:B:32:LEU:HD13	1:B:61:LEU:HD21	1.98	0.46
1:A:48:LYS:O	1:A:49:TYR:HB2	2.16	0.46
1:J:75:TRP:CZ3	1:J:190:ILE:HD11	2.50	0.46
1:I:172:PRO:HB3	1:I:174:PHE:CZ	2.50	0.46
1:H:179:GLU:CG	1:H:180:GLU:N	2.79	0.46
1:L:136:THR:O	1:L:139:VAL:HG13	2.16	0.46
1:H:21:ALA:HB1	1:H:69:VAL:HG13	1.98	0.46
1:J:90:ASN:CG	1:J:93:SER:HB2	2.36	0.46
1:F:90:ASN:CG	1:F:93:SER:HB2	2.37	0.46
1:I:33:TYR:HD1	1:I:43:PHE:CD2	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ARG:HH12	1:A:133:ARG:HH22	1.64	0.46
1:K:129:ARG:O	1:K:133:ARG:HB2	2.16	0.46
1:L:58:ASP:O	1:L:62:ILE:HG13	2.16	0.46
1:H:64:TYR:CE1	1:H:172:PRO:O	2.68	0.46
1:B:63:LYS:HE3	1:H:63:LYS:HZ3	1.81	0.46
1:C:188:THR:C	1:C:189:THR:HG23	2.37	0.46
1:L:163:VAL:O	1:L:164:PRO:C	2.54	0.46
1:F:58:ASP:O	1:F:62:ILE:HG13	2.16	0.46
1:B:179:GLU:CG	1:B:180:GLU:N	2.79	0.46
1:B:180:GLU:HA	1:B:180:GLU:OE1	2.16	0.46
1:H:75:TRP:CD1	1:H:161:LEU:HD21	2.51	0.45
1:K:153:LEU:O	1:K:154:LEU:HD23	2.17	0.45
1:C:42:THR:HB	1:C:43:PHE:CE1	2.52	0.45
1:J:32:LEU:HD13	1:J:61:LEU:HD21	1.98	0.45
1:I:131:VAL:HG12	1:I:132:ILE:N	2.31	0.45
1:G:11:ILE:HD13	1:G:11:ILE:HA	1.77	0.45
1:I:34:GLN:HE22	1:I:136:THR:HA	1.81	0.45
1:K:38:TYR:CD1	1:K:61:LEU:HD22	2.51	0.45
1:G:56:THR:HG21	1:G:61:LEU:CD2	2.41	0.45
1:G:27:GLY:O	1:G:31:ILE:HG13	2.16	0.45
1:D:136:THR:O	1:D:139:VAL:HG13	2.15	0.45
1:L:167:TRP:CH2	1:L:172:PRO:HD2	2.52	0.45
1:L:64:TYR:HE1	1:L:172:PRO:O	1.98	0.45
1:D:90:ASN:CG	1:D:93:SER:CB	2.85	0.45
1:J:14:ARG:HB3	1:J:105:GLU:CB	2.45	0.45
1:B:93:SER:OG	1:B:95:GLU:HG2	2.17	0.45
1:B:33:TYR:CD2	1:B:54:LEU:HD23	2.51	0.45
1:C:129:ARG:O	1:C:133:ARG:HB2	2.17	0.45
1:L:75:TRP:CZ3	1:L:190:ILE:HD11	2.51	0.45
1:G:56:THR:CG2	1:G:61:LEU:HD23	2.42	0.45
1:J:58:ASP:O	1:J:62:ILE:HG13	2.17	0.45
1:L:59:LEU:HA	1:L:59:LEU:HD12	1.84	0.45
1:G:34:GLN:OE1	1:G:136:THR:HA	2.16	0.45
1:C:178:SER:O	1:C:201:ILE:HD11	2.16	0.45
1:K:56:THR:CG2	1:K:61:LEU:HD23	2.41	0.45
1:F:120:SER:HB3	1:I:58:ASP:OD1	2.16	0.45
1:I:28:ILE:HD13	1:I:153:LEU:HD21	1.97	0.45
1:E:34:GLN:OE1	1:E:136:THR:HA	2.16	0.45
1:B:160:ASP:O	1:B:161:LEU:HG	2.16	0.45
1:L:78:LYS:HG3	1:L:80:SER:OG	2.17	0.45
1:G:135:ILE:O	1:G:138:THR:HG22	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:139:VAL:CG1	1:I:140:THR:N	2.80	0.45
1:D:19:ILE:HG23	1:D:20:VAL:N	2.31	0.45
1:A:54:LEU:N	1:A:54:LEU:HD12	2.32	0.45
1:I:20:VAL:O	1:I:24:PHE:HD1	1.99	0.45
1:K:24:PHE:CD2	1:K:84:LEU:HG	2.52	0.45
1:D:161:LEU:C	1:D:163:VAL:H	2.19	0.45
1:F:72:LEU:HD21	1:F:84:LEU:HD11	1.99	0.45
1:K:54:LEU:N	1:K:54:LEU:HD12	2.31	0.45
1:K:34:GLN:OE1	1:K:136:THR:HA	2.16	0.45
1:G:131:VAL:HG12	1:G:132:ILE:N	2.31	0.45
1:I:142:LEU:O	1:I:143:PRO:O	2.34	0.45
1:H:75:TRP:CZ3	1:H:190:ILE:HD11	2.51	0.45
1:H:90:ASN:CB	1:H:93:SER:HB3	2.39	0.45
1:C:49:TYR:O	1:C:129:ARG:CD	2.61	0.45
1:A:56:THR:CG2	1:A:61:LEU:HD23	2.42	0.45
1:D:159:LYS:HA	1:D:192:LYS:HB3	1.99	0.45
1:G:64:TYR:C	1:G:64:TYR:HD1	2.20	0.45
1:L:39:PRO:HD3	1:L:181:VAL:HG22	1.99	0.45
1:E:131:VAL:HG12	1:E:132:ILE:N	2.32	0.45
1:F:137:ALA:O	1:F:140:THR:HB	2.17	0.45
1:L:167:TRP:O	1:L:167:TRP:CE3	2.70	0.45
1:F:90:ASN:CB	1:F:93:SER:HB3	2.39	0.45
1:A:100:TRP:CE3	1:A:197:VAL:HG23	2.52	0.45
1:E:28:ILE:HD13	1:E:153:LEU:HD21	1.99	0.45
1:I:175:ILE:HB	1:I:178:SER:HB3	1.98	0.45
1:B:136:THR:O	1:B:139:VAL:HG13	2.16	0.45
1:L:138:THR:C	1:L:140:THR:N	2.70	0.45
1:B:47:GLN:HB2	1:B:51:LEU:O	2.17	0.45
1:L:93:SER:OG	1:L:95:GLU:HG2	2.17	0.44
1:D:31:ILE:HD13	1:D:100:TRP:CE3	2.52	0.44
1:B:137:ALA:O	1:B:140:THR:HB	2.16	0.44
1:H:180:GLU:HA	1:H:180:GLU:OE1	2.17	0.44
1:C:175:ILE:HB	1:C:178:SER:HB3	1.97	0.44
1:D:21:ALA:HB1	1:D:69:VAL:HG13	1.98	0.44
1:J:21:ALA:HB1	1:J:69:VAL:HG13	1.99	0.44
1:D:48:LYS:O	1:D:49:TYR:HB2	2.17	0.44
1:G:117:ARG:HA	1:G:117:ARG:HD2	1.45	0.44
1:B:76:LEU:O	1:B:78:LYS:N	2.51	0.44
1:C:84:LEU:HD13	1:C:85:VAL:N	2.32	0.44
1:H:160:ASP:OD2	1:I:159:LYS:NZ	2.50	0.44
1:I:35:ARG:HH11	1:I:35:ARG:CB	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:THR:C	1:B:140:THR:N	2.70	0.44
1:A:32:LEU:HD22	1:A:37:ILE:CG2	2.39	0.44
1:E:32:LEU:HD22	1:E:37:ILE:CG2	2.39	0.44
1:L:72:LEU:HD21	1:L:84:LEU:HD11	1.99	0.44
1:D:184:ARG:HA	1:D:184:ARG:HD3	1.81	0.44
1:F:48:LYS:HA	1:F:48:LYS:HD2	1.79	0.44
1:J:90:ASN:ND2	1:J:93:SER:HB2	2.33	0.44
1:J:39:PRO:HD3	1:J:181:VAL:HG22	1.99	0.44
1:C:56:THR:HG21	1:C:61:LEU:CD2	2.46	0.44
1:B:91:ILE:CD1	1:B:150:SER:HB3	2.43	0.44
1:B:199:TYR:O	1:C:154:LEU:HD22	2.18	0.44
1:K:129:ARG:NH2	2:K:206:SO4:O2	2.49	0.44
1:F:75:TRP:CZ3	1:F:190:ILE:HD11	2.52	0.44
1:D:76:LEU:O	1:D:78:LYS:N	2.50	0.44
1:E:84:LEU:HD13	1:E:85:VAL:N	2.33	0.44
1:K:64:TYR:C	1:K:64:TYR:HD1	2.21	0.44
1:L:197:VAL:HG11	1:L:199:TYR:CZ	2.53	0.44
1:L:28:ILE:O	1:L:32:LEU:HG	2.17	0.44
1:L:138:THR:O	1:L:140:THR:N	2.51	0.44
1:J:137:ALA:O	1:J:140:THR:HB	2.16	0.44
1:A:161:LEU:HG	1:A:162:VAL:N	2.33	0.44
1:B:59:LEU:HD12	1:B:59:LEU:HA	1.88	0.44
1:B:16:SER:O	1:B:19:ILE:HG22	2.18	0.44
1:J:93:SER:OG	1:J:95:GLU:HG2	2.18	0.44
1:B:90:ASN:ND2	1:B:93:SER:HB2	2.33	0.44
1:H:163:VAL:O	1:H:164:PRO:C	2.55	0.44
1:K:133:ARG:NH2	2:K:206:SO4:O4	2.51	0.44
1:G:153:LEU:O	1:G:154:LEU:HD23	2.18	0.44
1:G:84:LEU:HD13	1:G:85:VAL:N	2.33	0.44
1:G:172:PRO:HB3	1:G:174:PHE:CZ	2.53	0.44
1:D:135:ILE:O	1:D:139:VAL:HG12	2.17	0.44
1:F:21:ALA:HB1	1:F:69:VAL:HG13	2.00	0.44
1:C:48:LYS:O	1:C:49:TYR:HB2	2.17	0.43
1:C:34:GLN:OE1	1:C:136:THR:HA	2.18	0.43
1:A:135:ILE:O	1:A:138:THR:HG22	2.17	0.43
1:G:171:GLY:HA2	1:G:172:PRO:HD3	1.83	0.43
1:J:161:LEU:C	1:J:163:VAL:H	2.20	0.43
1:B:21:ALA:HB1	1:B:69:VAL:HG13	2.00	0.43
1:B:65:LEU:HD23	1:B:65:LEU:HA	1.76	0.43
1:H:72:LEU:HD21	1:H:84:LEU:HD11	2.00	0.43
1:B:72:LEU:HD21	1:B:84:LEU:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:90:ASN:CG	1:H:93:SER:HB2	2.38	0.43
1:H:90:ASN:ND2	1:H:93:SER:HB2	2.34	0.43
1:A:28:ILE:HD13	1:A:153:LEU:HD21	2.00	0.43
1:D:160:ASP:O	1:D:161:LEU:HG	2.18	0.43
1:J:72:LEU:HD21	1:J:84:LEU:HD11	1.99	0.43
1:G:32:LEU:HD22	1:G:37:ILE:CG2	2.40	0.43
1:L:90:ASN:CG	1:L:93:SER:HB2	2.38	0.43
1:A:129:ARG:NH2	2:A:207:SO4:O2	2.51	0.43
1:A:49:TYR:O	1:A:129:ARG:CD	2.63	0.43
1:A:38:TYR:CD1	1:A:61:LEU:HD22	2.53	0.43
1:B:58:ASP:O	1:B:62:ILE:HG13	2.19	0.43
1:B:76:LEU:C	1:B:78:LYS:H	2.21	0.43
1:A:34:GLN:OE1	1:A:136:THR:HA	2.18	0.43
1:K:71:GLN:HG2	1:K:167:TRP:CZ2	2.54	0.43
1:J:160:ASP:O	1:J:161:LEU:HG	2.19	0.43
1:F:161:LEU:C	1:F:163:VAL:H	2.21	0.43
1:L:140:THR:HG22	1:L:141:PHE:CD1	2.53	0.43
1:F:38:TYR:CZ	1:F:61:LEU:HD22	2.54	0.43
1:K:27:GLY:O	1:K:31:ILE:HG13	2.17	0.43
1:K:24:PHE:O	1:K:28:ILE:HG13	2.18	0.43
1:E:153:LEU:O	1:E:154:LEU:HD23	2.17	0.43
1:J:24:PHE:O	1:J:28:ILE:HG22	2.19	0.43
1:J:76:LEU:C	1:J:78:LYS:H	2.21	0.43
1:C:117:ARG:HA	1:C:117:ARG:HD2	1.48	0.43
1:L:180:GLU:HA	1:L:180:GLU:OE1	2.19	0.43
1:E:64:TYR:C	1:E:64:TYR:HD1	2.22	0.43
1:H:28:ILE:O	1:H:32:LEU:HG	2.18	0.43
1:H:140:THR:HG22	1:H:141:PHE:CD1	2.54	0.43
1:J:121:GLN:HE21	1:J:125:GLN:HE21	1.66	0.43
1:A:108:LYS:HD2	1:A:111:LYS:HZ1	1.83	0.43
1:I:133:ARG:HD2	1:J:34:GLN:HA	1.98	0.43
1:K:172:PRO:CB	1:K:174:PHE:CZ	3.02	0.43
1:J:179:GLU:CG	1:J:180:GLU:H	2.32	0.43
1:F:180:GLU:OE1	1:F:180:GLU:HA	2.19	0.43
1:D:175:ILE:CG2	1:D:176:THR:N	2.82	0.43
1:L:90:ASN:ND2	1:L:93:SER:HB2	2.34	0.43
1:E:129:ARG:NH2	2:E:206:SO4:O3	2.49	0.43
1:A:59:LEU:CD2	1:E:44:THR:CG2	2.97	0.43
1:F:140:THR:HG22	1:F:141:PHE:CD1	2.54	0.43
1:B:175:ILE:HD13	1:B:175:ILE:HA	1.67	0.43
1:J:33:TYR:HB2	1:J:43:PHE:CD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:93:SER:OG	1:H:95:GLU:HG2	2.18	0.43
1:D:76:LEU:C	1:D:78:LYS:H	2.21	0.43
1:J:159:LYS:HA	1:J:192:LYS:HB3	2.01	0.43
1:L:179:GLU:CG	1:L:180:GLU:H	2.32	0.43
1:D:138:THR:C	1:D:140:THR:N	2.71	0.43
1:J:47:GLN:HB2	1:J:51:LEU:O	2.19	0.43
1:C:71:GLN:HG2	1:C:167:TRP:CZ2	2.54	0.42
1:I:171:GLY:HA2	1:I:172:PRO:HD3	1.84	0.42
1:L:31:ILE:HD13	1:L:100:TRP:CE3	2.54	0.42
1:F:138:THR:C	1:F:140:THR:N	2.72	0.42
1:L:71:GLN:HG2	1:L:167:TRP:CD1	2.54	0.42
1:H:16:SER:O	1:H:20:VAL:HG23	2.19	0.42
1:H:31:ILE:HD13	1:H:100:TRP:CE3	2.54	0.42
1:F:179:GLU:CG	1:F:180:GLU:H	2.32	0.42
1:L:76:LEU:C	1:L:78:LYS:N	2.73	0.42
1:J:76:LEU:O	1:J:78:LYS:N	2.52	0.42
1:L:173:GLN:HB2	1:L:187:THR:HG22	2.00	0.42
1:G:129:ARG:HH12	1:G:133:ARG:HH22	1.65	0.42
1:C:56:THR:CG2	1:C:61:LEU:HD23	2.46	0.42
1:A:27:GLY:O	1:A:31:ILE:HG13	2.18	0.42
1:D:179:GLU:CG	1:D:180:GLU:H	2.31	0.42
1:I:183:LEU:HB2	1:I:195:SER:O	2.18	0.42
1:G:48:LYS:HD2	1:G:48:LYS:HA	1.90	0.42
1:L:157:THR:HG21	1:L:161:LEU:HD23	2.01	0.42
1:K:10:SER:CA	1:K:117:ARG:O	2.64	0.42
1:A:59:LEU:CD2	1:E:44:THR:HG21	2.49	0.42
1:C:153:LEU:O	1:C:154:LEU:HD23	2.20	0.42
1:D:121:GLN:HE21	1:D:125:GLN:HE21	1.66	0.42
1:E:134:GLN:HE21	1:F:139:VAL:HG13	1.83	0.42
1:J:28:ILE:O	1:J:32:LEU:HG	2.20	0.42
1:I:18:GLU:HG2	1:I:73:LYS:HD2	1.99	0.42
1:F:93:SER:OG	1:F:95:GLU:HG2	2.19	0.42
1:A:172:PRO:HB3	1:A:174:PHE:CZ	2.54	0.42
1:K:84:LEU:HD13	1:K:84:LEU:C	2.39	0.42
1:C:108:LYS:HD2	1:C:111:LYS:HZ1	1.85	0.42
1:F:179:GLU:HG3	1:F:180:GLU:N	2.35	0.42
1:D:175:ILE:HG22	1:D:176:THR:N	2.33	0.42
1:L:184:ARG:HD3	1:L:184:ARG:HA	1.85	0.42
1:I:34:GLN:NE2	1:I:139:VAL:HB	2.35	0.42
1:B:90:ASN:CG	1:B:93:SER:CB	2.88	0.42
1:D:58:ASP:O	1:D:62:ILE:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:63:LYS:HG3	1:L:59:LEU:CD2	2.46	0.42
1:B:76:LEU:C	1:B:78:LYS:N	2.73	0.42
1:A:84:LEU:HD13	1:A:85:VAL:N	2.34	0.42
1:B:27:GLY:O	1:B:31:ILE:HG13	2.19	0.42
1:F:59:LEU:HD12	1:F:59:LEU:HA	1.88	0.42
1:L:48:LYS:O	1:L:49:TYR:HB2	2.20	0.42
1:D:188:THR:HG22	1:D:189:THR:N	2.35	0.42
1:C:172:PRO:HB3	1:C:174:PHE:CZ	2.54	0.42
1:D:77:TYR:HE2	1:D:78:LYS:HE2	1.85	0.42
1:F:32:LEU:HD13	1:F:61:LEU:HD21	2.01	0.42
1:L:47:GLN:HB2	1:L:51:LEU:O	2.19	0.42
1:I:35:ARG:CB	1:I:35:ARG:CZ	2.87	0.42
1:G:49:TYR:O	1:G:129:ARG:CD	2.62	0.42
1:G:71:GLN:HG2	1:G:167:TRP:CZ2	2.55	0.42
1:D:179:GLU:HG3	1:D:180:GLU:N	2.35	0.42
1:H:138:THR:C	1:H:140:THR:N	2.72	0.42
1:L:76:LEU:C	1:L:78:LYS:H	2.22	0.42
1:B:138:THR:O	1:B:140:THR:N	2.53	0.42
1:B:203:VAL:O	1:B:203:VAL:HG12	2.20	0.42
1:I:11:ILE:HA	1:I:11:ILE:HD13	1.78	0.42
1:L:16:SER:O	1:L:20:VAL:HG23	2.19	0.42
1:B:39:PRO:HD2	1:B:42:THR:OG1	2.19	0.42
1:E:26:PHE:HD1	1:E:49:TYR:CE1	2.38	0.42
1:C:10:SER:CA	1:C:117:ARG:O	2.63	0.42
1:I:117:ARG:HA	1:I:117:ARG:HD2	1.44	0.42
1:B:31:ILE:HD13	1:B:100:TRP:CE3	2.55	0.42
1:J:179:GLU:HG3	1:J:180:GLU:N	2.35	0.42
1:J:138:THR:C	1:J:140:THR:N	2.73	0.42
1:B:28:ILE:O	1:B:32:LEU:HG	2.20	0.42
1:E:11:ILE:HA	1:E:11:ILE:HD13	1.78	0.42
1:I:49:TYR:O	1:I:129:ARG:CD	2.62	0.42
1:B:188:THR:HG22	1:B:189:THR:N	2.35	0.42
1:D:76:LEU:C	1:D:78:LYS:N	2.72	0.42
1:I:84:LEU:HD13	1:I:85:VAL:N	2.35	0.42
1:C:84:LEU:C	1:C:84:LEU:HD13	2.40	0.42
1:E:71:GLN:HG2	1:E:167:TRP:CZ2	2.55	0.42
1:B:161:LEU:C	1:B:163:VAL:H	2.22	0.42
1:F:186:PHE:CD2	1:F:186:PHE:C	2.94	0.41
1:I:86:VAL:HB	1:I:100:TRP:HB2	2.01	0.41
1:J:180:GLU:HA	1:J:180:GLU:OE1	2.19	0.41
1:L:21:ALA:HB1	1:L:69:VAL:HG13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:78:LYS:HG3	1:F:80:SER:OG	2.19	0.41
1:D:203:VAL:O	1:D:203:VAL:HG12	2.20	0.41
1:L:179:GLU:HG3	1:L:180:GLU:N	2.35	0.41
1:J:31:ILE:HD13	1:J:100:TRP:CE3	2.55	0.41
1:D:38:TYR:CE1	1:D:61:LEU:HD22	2.55	0.41
1:G:18:GLU:HG2	1:G:73:LYS:HD2	2.01	0.41
1:F:184:ARG:HA	1:F:184:ARG:HD3	1.84	0.41
1:F:45:ARG:O	1:I:146:GLU:OE2	2.38	0.41
1:I:39:PRO:HD2	1:I:42:THR:OG1	2.19	0.41
1:A:48:LYS:HA	1:A:48:LYS:HD2	1.88	0.41
1:H:76:LEU:C	1:H:78:LYS:H	2.23	0.41
1:C:84:LEU:CD1	1:C:84:LEU:C	2.89	0.41
1:B:140:THR:HG22	1:B:141:PHE:CD1	2.55	0.41
1:F:76:LEU:C	1:F:78:LYS:N	2.74	0.41
1:K:11:ILE:HA	1:K:11:ILE:HD13	1.75	0.41
1:J:132:ILE:HD13	1:J:132:ILE:N	2.35	0.41
1:D:173:GLN:CB	1:D:187:THR:HG22	2.42	0.41
1:D:14:ARG:HB3	1:D:105:GLU:HB3	2.02	0.41
1:D:186:PHE:C	1:D:186:PHE:CD2	2.94	0.41
1:L:76:LEU:O	1:L:78:LYS:N	2.53	0.41
1:K:135:ILE:O	1:K:138:THR:HG22	2.20	0.41
1:F:186:PHE:HD2	1:F:187:THR:N	2.19	0.41
1:B:77:TYR:HE2	1:B:78:LYS:HE2	1.86	0.41
1:H:14:ARG:HB3	1:H:105:GLU:HB3	2.01	0.41
1:A:64:TYR:C	1:A:64:TYR:HD1	2.24	0.41
1:H:77:TYR:HE2	1:H:78:LYS:HE2	1.86	0.41
1:L:14:ARG:HB3	1:L:105:GLU:CG	2.49	0.41
1:B:179:GLU:CG	1:B:180:GLU:H	2.33	0.41
1:F:121:GLN:HE21	1:F:125:GLN:HE21	1.68	0.41
1:B:132:ILE:N	1:B:132:ILE:HD13	2.35	0.41
1:J:16:SER:HB2	1:J:109:THR:CG2	2.36	0.41
1:F:14:ARG:HB3	1:F:105:GLU:HB3	2.03	0.41
1:B:163:VAL:HA	1:B:164:PRO:HD3	1.94	0.41
1:F:24:PHE:O	1:F:28:ILE:HG22	2.20	0.41
1:J:76:LEU:C	1:J:78:LYS:N	2.73	0.41
1:L:173:GLN:CB	1:L:187:THR:HG22	2.50	0.41
1:H:184:ARG:HA	1:H:184:ARG:HD3	1.85	0.41
1:I:37:ILE:HG13	1:I:37:ILE:O	2.20	0.41
1:F:45:ARG:O	1:I:146:GLU:CD	2.59	0.41
1:J:39:PRO:HD2	1:J:42:THR:OG1	2.20	0.41
1:C:51:LEU:HB2	1:C:53:LEU:HD21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:39:PRO:HD2	1:D:42:THR:OG1	2.21	0.41
1:F:159:LYS:HA	1:F:192:LYS:HB3	2.02	0.41
1:H:179:GLU:CG	1:H:180:GLU:H	2.33	0.41
1:B:38:TYR:CZ	1:B:61:LEU:HD22	2.56	0.41
1:F:76:LEU:C	1:F:78:LYS:H	2.23	0.41
1:L:201:ILE:O	1:L:201:ILE:HG23	2.21	0.41
1:F:48:LYS:HE3	1:I:37:ILE:O	2.21	0.41
1:I:141:PHE:O	1:J:129:ARG:CD	2.65	0.41
1:B:200:LYS:HG3	1:C:170:SER:CB	2.51	0.41
1:I:33:TYR:CE2	1:I:54:LEU:HD13	2.55	0.41
1:A:188:THR:C	1:A:189:THR:HG23	2.40	0.41
1:K:49:TYR:O	1:K:129:ARG:CD	2.63	0.41
1:H:39:PRO:HD2	1:H:42:THR:OG1	2.21	0.41
1:C:171:GLY:HA2	1:C:172:PRO:HD3	1.81	0.41
1:L:14:ARG:HB3	1:L:105:GLU:HB3	2.03	0.41
1:K:84:LEU:C	1:K:84:LEU:CD1	2.89	0.41
1:F:28:ILE:O	1:F:32:LEU:HG	2.20	0.41
1:D:175:ILE:HA	1:D:175:ILE:HD13	1.68	0.41
1:F:76:LEU:O	1:F:78:LYS:N	2.54	0.41
1:A:12:THR:OG1	1:A:14:ARG:HB3	2.21	0.41
1:A:131:VAL:HG12	1:A:132:ILE:N	2.34	0.41
1:F:197:VAL:HG12	1:F:199:TYR:CE2	2.56	0.41
1:E:140:THR:OG1	1:F:53:LEU:HD23	2.21	0.41
1:K:95:GLU:HG2	1:K:96:VAL:N	2.34	0.41
1:G:76:LEU:HD23	1:G:76:LEU:HA	1.89	0.41
1:D:39:PRO:HD3	1:D:181:VAL:HG22	2.02	0.41
1:K:129:ARG:HH12	1:K:133:ARG:HH22	1.67	0.41
1:K:48:LYS:O	1:K:49:TYR:HB2	2.21	0.41
1:J:14:ARG:HB3	1:J:105:GLU:HB3	2.03	0.41
1:F:138:THR:O	1:F:140:THR:N	2.54	0.41
1:B:194:ASN:HB3	1:B:195:SER:H	1.77	0.41
1:B:104:ILE:HG21	1:B:104:ILE:HD13	1.83	0.41
1:I:54:LEU:N	1:I:54:LEU:HD12	2.36	0.40
1:F:31:ILE:HD13	1:F:100:TRP:CE3	2.55	0.40
1:E:18:GLU:HG2	1:E:73:LYS:HD2	2.01	0.40
1:K:18:GLU:HG2	1:K:73:LYS:HD2	2.03	0.40
1:H:76:LEU:C	1:H:78:LYS:N	2.74	0.40
1:G:161:LEU:HG	1:G:162:VAL:N	2.36	0.40
1:D:72:LEU:HD21	1:D:84:LEU:HD11	2.02	0.40
1:I:39:PRO:O	1:I:40:SER:C	2.60	0.40
1:F:186:PHE:HD2	1:F:186:PHE:C	2.24	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:199:TYR:O	1:I:154:LEU:HA	2.22	0.40
1:K:64:TYR:HE1	1:K:68:VAL:HG21	1.86	0.40
1:C:76:LEU:HD23	1:C:76:LEU:HA	1.92	0.40
1:I:43:PHE:HB3	1:I:54:LEU:HB3	2.02	0.40
1:K:171:GLY:HA2	1:K:172:PRO:HD3	1.84	0.40
1:K:117:ARG:HA	1:K:117:ARG:HD2	1.44	0.40
1:D:121:GLN:HB3	1:K:43:PHE:O	2.21	0.40
1:J:161:LEU:HD23	1:J:161:LEU:HA	1.90	0.40
1:H:138:THR:O	1:H:140:THR:N	2.53	0.40
1:A:18:GLU:HG2	1:A:73:LYS:HD2	2.02	0.40
1:C:103:ASP:HB2	1:C:194:ASN:HB2	2.04	0.40
1:K:161:LEU:HG	1:K:162:VAL:N	2.36	0.40
1:E:95:GLU:HG2	1:E:96:VAL:N	2.35	0.40
1:J:203:VAL:O	1:J:203:VAL:HG12	2.22	0.40
1:F:42:THR:HB	1:F:43:PHE:CE1	2.57	0.40
1:G:49:TYR:CD2	1:G:125:GLN:HG3	2.56	0.40
1:H:33:TYR:HB2	1:H:43:PHE:CD2	2.56	0.40
1:H:58:ASP:O	1:H:62:ILE:HG13	2.21	0.40
1:H:179:GLU:HG3	1:H:180:GLU:N	2.36	0.40
1:E:175:ILE:O	1:E:176:THR:C	2.60	0.40
1:C:64:TYR:C	1:C:64:TYR:HD1	2.24	0.40
1:H:132:ILE:HD13	1:H:132:ILE:N	2.35	0.40
1:K:76:LEU:HA	1:K:76:LEU:HD23	1.91	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	191/207 (92%)	173 (91%)	18 (9%)	0	100	100
1	B	181/207 (87%)	156 (86%)	25 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	193/207 (93%)	174 (90%)	18 (9%)	1 (0%)	34	76
1	D	183/207 (88%)	158 (86%)	25 (14%)	0	100	100
1	E	184/207 (89%)	167 (91%)	16 (9%)	1 (0%)	34	76
1	F	181/207 (87%)	159 (88%)	21 (12%)	1 (1%)	30	73
1	G	183/207 (88%)	169 (92%)	13 (7%)	1 (0%)	34	76
1	H	182/207 (88%)	154 (85%)	28 (15%)	0	100	100
1	I	184/207 (89%)	165 (90%)	16 (9%)	3 (2%)	12	57
1	J	181/207 (87%)	158 (87%)	22 (12%)	1 (1%)	30	73
1	K	184/207 (89%)	167 (91%)	16 (9%)	1 (0%)	34	76
1	L	183/207 (88%)	157 (86%)	26 (14%)	0	100	100
All	All	2210/2484 (89%)	1957 (89%)	244 (11%)	9 (0%)	39	79

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	176	THR
1	G	176	THR
1	I	176	THR
1	J	169	GLU
1	K	176	THR
1	C	176	THR
1	F	169	GLU
1	I	143	PRO
1	I	37	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	180/192 (94%)	172 (96%)	8 (4%)	35	72
1	B	173/192 (90%)	161 (93%)	12 (7%)	19	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	182/192 (95%)	172 (94%)	10 (6%)	27	66
1	D	175/192 (91%)	163 (93%)	12 (7%)	19	59
1	E	176/192 (92%)	170 (97%)	6 (3%)	44	77
1	F	173/192 (90%)	159 (92%)	14 (8%)	15	53
1	G	175/192 (91%)	168 (96%)	7 (4%)	38	73
1	H	174/192 (91%)	164 (94%)	10 (6%)	25	65
1	I	176/192 (92%)	170 (97%)	6 (3%)	44	77
1	J	173/192 (90%)	163 (94%)	10 (6%)	25	64
1	K	176/192 (92%)	170 (97%)	6 (3%)	44	77
1	L	174/192 (91%)	164 (94%)	10 (6%)	25	65
All	All	2107/2304 (91%)	1996 (95%)	111 (5%)	28	67

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	40	SER
1	A	64	TYR
1	A	84	LEU
1	A	117	ARG
1	A	138	THR
1	A	144	LEU
1	A	167	TRP
1	B	14	ARG
1	B	43	PHE
1	B	66	ASN
1	B	105	GLU
1	B	130	SER
1	B	139	VAL
1	B	151	PHE
1	B	153	LEU
1	B	165	GLU
1	B	166	LYS
1	B	185	SER
1	B	186	PHE
1	C	11	ILE
1	C	40	SER
1	C	45	ARG

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Mol	Chain	Res	Type
1	C	53	LEU
1	C	64	TYR
1	C	84	LEU
1	C	117	ARG
1	C	138	THR
1	C	144	LEU
1	C	167	TRP
1	D	14	ARG
1	D	43	PHE
1	D	66	ASN
1	D	105	GLU
1	D	130	SER
1	D	139	VAL
1	D	151	PHE
1	D	153	LEU
1	D	165	GLU
1	D	166	LYS
1	D	185	SER
1	D	186	PHE
1	E	40	SER
1	E	64	TYR
1	E	84	LEU
1	E	138	THR
1	E	144	LEU
1	E	167	TRP
1	F	14	ARG
1	F	43	PHE
1	F	47	GLN
1	F	48	LYS
1	F	49	TYR
1	F	66	ASN
1	F	105	GLU
1	F	130	SER
1	F	139	VAL
1	F	151	PHE
1	F	153	LEU
1	F	165	GLU
1	F	166	LYS
1	F	186	PHE
1	G	40	SER
1	G	64	TYR
1	G	84	LEU

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Mol	Chain	Res	Type
1	G	117	ARG
1	G	138	THR
1	G	144	LEU
1	G	167	TRP
1	H	14	ARG
1	H	30	SER
1	H	43	PHE
1	H	105	GLU
1	H	130	SER
1	H	139	VAL
1	H	151	PHE
1	H	153	LEU
1	H	166	LYS
1	H	186	PHE
1	I	35	ARG
1	I	64	TYR
1	I	84	LEU
1	I	117	ARG
1	I	144	LEU
1	I	167	TRP
1	J	14	ARG
1	J	43	PHE
1	J	66	ASN
1	J	105	GLU
1	J	139	VAL
1	J	151	PHE
1	J	153	LEU
1	J	165	GLU
1	J	166	LYS
1	J	186	PHE
1	K	40	SER
1	K	64	TYR
1	K	84	LEU
1	K	117	ARG
1	K	138	THR
1	K	144	LEU
1	L	14	ARG
1	L	43	PHE
1	L	66	ASN
1	L	105	GLU
1	L	130	SER
1	L	139	VAL

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Mol	Chain	Res	Type
1	L	151	PHE
1	L	153	LEU
1	L	166	LYS
1	L	186	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	125	GLN
1	D	125	GLN
1	F	121	GLN
1	J	121	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	206	-	4,4,4	0.19	0	6,6,6	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	207	-	4,4,4	0.33	0	6,6,6	0.31	0
2	SO4	B	206	-	4,4,4	0.19	0	6,6,6	0.31	0
2	SO4	C	206	-	4,4,4	0.32	0	6,6,6	0.40	0
2	SO4	C	207	-	4,4,4	0.18	0	6,6,6	0.14	0
2	SO4	D	206	-	4,4,4	0.22	0	6,6,6	0.50	0
2	SO4	E	206	-	4,4,4	0.27	0	6,6,6	0.16	0
2	SO4	G	206	-	4,4,4	0.27	0	6,6,6	0.32	0
2	SO4	H	206	-	4,4,4	0.16	0	6,6,6	0.40	0
2	SO4	I	206	-	4,4,4	0.25	0	6,6,6	0.15	0
2	SO4	K	206	-	4,4,4	0.24	0	6,6,6	0.22	0
2	SO4	L	206	-	4,4,4	0.19	0	6,6,6	0.34	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	206	-	-	0/0/0/0	0/0/0/0
2	SO4	A	207	-	-	0/0/0/0	0/0/0/0
2	SO4	B	206	-	-	0/0/0/0	0/0/0/0
2	SO4	C	206	-	-	0/0/0/0	0/0/0/0
2	SO4	C	207	-	-	0/0/0/0	0/0/0/0
2	SO4	D	206	-	-	0/0/0/0	0/0/0/0
2	SO4	E	206	-	-	0/0/0/0	0/0/0/0
2	SO4	G	206	-	-	0/0/0/0	0/0/0/0
2	SO4	H	206	-	-	0/0/0/0	0/0/0/0
2	SO4	I	206	-	-	0/0/0/0	0/0/0/0
2	SO4	K	206	-	-	0/0/0/0	0/0/0/0
2	SO4	L	206	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	207	SO4	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	206	SO4	2	0
2	E	206	SO4	2	0
2	G	206	SO4	1	0
2	K	206	SO4	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	193/207 (93%)	-0.05	4 (2%) 67 55	71, 120, 205, 262	0
1	B	185/207 (89%)	-0.04	3 (1%) 74 63	69, 104, 175, 238	0
1	C	195/207 (94%)	0.17	4 (2%) 67 55	69, 129, 220, 348	0
1	D	187/207 (90%)	-0.01	4 (2%) 67 55	67, 101, 187, 271	0
1	E	188/207 (90%)	0.51	18 (9%) 10 7	104, 186, 265, 315	0
1	F	185/207 (89%)	0.38	20 (10%) 8 6	96, 149, 241, 315	0
1	G	187/207 (90%)	0.35	10 (5%) 30 22	104, 175, 247, 312	0
1	H	186/207 (89%)	0.41	16 (8%) 13 10	92, 159, 236, 277	0
1	I	188/207 (90%)	0.59	20 (10%) 8 6	111, 180, 273, 324	0
1	J	185/207 (89%)	0.49	24 (12%) 5 5	108, 169, 246, 283	0
1	K	188/207 (90%)	0.36	13 (6%) 20 13	94, 176, 250, 315	0
1	L	187/207 (90%)	0.41	16 (8%) 13 10	81, 146, 241, 396	0
All	All	2254/2484 (90%)	0.30	152 (6%) 21 14	67, 151, 245, 396	0

All (152) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	204	ASN	8.1
1	F	9	GLY	5.7
1	F	109	THR	5.4
1	E	169	GLU	4.6
1	D	109	THR	4.5
1	J	162	VAL	4.5
1	I	118	GLU	4.5
1	I	110	ALA	4.3
1	L	9	GLY	4.2
1	I	109	THR	4.1
1	J	200	LYS	4.0

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Mol	Chain	Res	Type	RSRZ
1	E	165	GLU	4.0
1	K	119	LYS	4.0
1	I	203	VAL	4.0
1	J	9	GLY	4.0
1	B	9	GLY	3.9
1	F	177	ASN	3.9
1	L	77	TYR	3.9
1	L	109	THR	3.8
1	L	199	TYR	3.6
1	L	157	THR	3.6
1	G	97	LEU	3.6
1	L	177	ASN	3.6
1	I	107	ASP	3.6
1	E	159	LYS	3.6
1	F	180	GLU	3.5
1	E	44	THR	3.4
1	J	161	LEU	3.4
1	K	169	GLU	3.4
1	F	195	SER	3.4
1	F	148	SER	3.3
1	F	184	ARG	3.3
1	L	161	LEU	3.3
1	E	56	THR	3.3
1	H	160	ASP	3.2
1	I	121	GLN	3.2
1	D	77	TYR	3.2
1	K	168	GLU	3.1
1	G	153	LEU	3.1
1	I	191	HIS	3.1
1	J	177	ASN	3.0
1	L	191	HIS	3.0
1	F	199	TYR	3.0
1	K	96	VAL	3.0
1	C	204	ASN	3.0
1	G	96	VAL	2.9
1	F	185	SER	2.9
1	J	175	ILE	2.9
1	E	109	THR	2.9
1	C	112	ASP	2.9
1	F	200	LYS	2.8
1	J	160	ASP	2.8
1	J	203	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	H	177	ASN	2.8
1	I	119	LYS	2.8
1	J	201	ILE	2.8
1	K	10	SER	2.7
1	A	116	PRO	2.7
1	G	200	LYS	2.7
1	F	176	THR	2.7
1	H	8	HIS	2.7
1	I	167	TRP	2.7
1	J	174	PHE	2.7
1	K	166	LYS	2.6
1	E	38	TYR	2.6
1	J	202	PRO	2.6
1	J	158	ASP	2.6
1	F	77	TYR	2.6
1	J	185	SER	2.6
1	F	169	GLU	2.6
1	E	167	TRP	2.6
1	L	192	LYS	2.6
1	F	202	PRO	2.6
1	H	179	GLU	2.6
1	I	54	LEU	2.6
1	L	190	ILE	2.6
1	J	184	ARG	2.5
1	I	11	ILE	2.5
1	J	91	ILE	2.5
1	H	200	LYS	2.5
1	H	190	ILE	2.5
1	E	118	GLU	2.5
1	F	92	GLU	2.5
1	J	109	THR	2.4
1	A	114	SER	2.4
1	J	195	SER	2.4
1	B	109	THR	2.4
1	E	12	THR	2.4
1	H	159	LYS	2.4
1	E	166	LYS	2.4
1	G	11	ILE	2.4
1	F	90	ASN	2.4
1	F	91	ILE	2.4
1	K	159	LYS	2.4
1	I	44	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	I	190	ILE	2.4
1	J	75	TRP	2.4
1	K	12	THR	2.4
1	K	118	GLU	2.4
1	H	161	LEU	2.3
1	L	91	ILE	2.3
1	H	194	ASN	2.3
1	J	163	VAL	2.3
1	H	191	HIS	2.3
1	E	168	GLU	2.3
1	I	159	LYS	2.3
1	I	49	TYR	2.3
1	K	165	GLU	2.3
1	K	97	LEU	2.3
1	I	168	GLU	2.3
1	J	18	GLU	2.3
1	E	188	THR	2.3
1	A	92	GLU	2.3
1	H	95	GLU	2.3
1	I	187	THR	2.3
1	F	203	VAL	2.3
1	H	165	GLU	2.3
1	H	9	GLY	2.3
1	E	158	ASP	2.2
1	J	169	GLU	2.2
1	L	92	GLU	2.2
1	H	192	LYS	2.2
1	L	90	ASN	2.2
1	B	77	TYR	2.2
1	G	152	ASP	2.2
1	C	10	SER	2.2
1	E	77	TYR	2.2
1	G	90	ASN	2.2
1	G	173	GLN	2.2
1	E	92	GLU	2.2
1	A	10	SER	2.1
1	F	201	ILE	2.1
1	G	149	CYS	2.1
1	H	97	LEU	2.1
1	D	108	LYS	2.1
1	I	106	CYS	2.1
1	C	203	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	202	PRO	2.1
1	K	117	ARG	2.1
1	E	124	ILE	2.1
1	L	179	GLU	2.1
1	I	13	LEU	2.1
1	L	110	ALA	2.1
1	L	80	SER	2.0
1	F	159	LYS	2.0
1	J	196	MET	2.0
1	H	82	GLN	2.0
1	J	165	GLU	2.0
1	K	13	LEU	2.0
1	I	10	SER	2.0
1	G	110	ALA	2.0
1	J	159	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	D	206	5/5	0.79	0.44	2.14	138,138,138,138	0
2	SO4	L	206	5/5	0.90	0.37	1.95	121,121,121,121	0
2	SO4	G	206	5/5	0.82	0.42	1.44	184,184,184,184	0
2	SO4	H	206	5/5	0.85	0.36	1.28	133,133,133,133	0
2	SO4	A	206	5/5	0.85	0.24	0.35	185,185,185,185	0
2	SO4	B	206	5/5	0.84	0.33	0.13	138,138,138,138	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	K	206	5/5	0.77	0.27	-0.05	196,196,196,196	0
2	SO4	E	206	5/5	0.68	0.26	-0.45	188,188,188,188	0
2	SO4	A	207	5/5	0.88	0.21	-0.59	128,128,128,128	0
2	SO4	C	206	5/5	0.88	0.20	-0.67	120,120,120,120	0
2	SO4	C	207	5/5	0.88	0.18	-0.69	191,191,191,191	0
2	SO4	I	206	5/5	0.80	0.18	-1.05	192,192,192,192	0

6.5 Other polymers [i](#)

There are no such residues in this entry.